
*
***** G A M M A S P E C T R U M A N A L Y S I S

*

OYSTER CREEK NUCLEAR GENERATING STATION CHEMISTRY DEPARTMENT

Detector Name : DET01
Report Generated On : 4/18/09 1:36:19 PM
Spectral Data File Name : C:\PCNT2K\CAMFILES\1LSOLMDA
 : *Co-60 1.13 E-6*
 : *Cs-137 1.46 E-7*
Sample Title : 1-L Solid Releas
Sample Description : Soil near CST; 2' depth
User ID : kg
Sample Type : 20
Sample Geometry : 1-L Marinelli
Peak Locate Threshold : 3.00
Peak Locate Range (in channels) : 1 - 4096
Peak Area Range (in channels) : 100 - 4096
Identification Energy Tolerance : 1.000 keV
Sample Size : 940.00 grams
Sample Taken On : 4/17/09 8:00:18 PM
Acquisition Started : 4/18/09 1:11:20 PM
Decay Time : 1.03E+003 Minutes
Live Time : 1796.0 Seconds
Real Time : 1797.1 Seconds
Dead Time : 0.06 %

Energy Calibration Used Done On : 1/8/09
Efficiency Calibration Used Done On : 1/8/09



**
***** B A C K G R O U N D S U B T R A C T R E P O R T

**

Detector Name: DET01
Sample Title: 1-L Solid Release
Peak Analysis Performed on: 4/18/09 1:36:19 PM

Peak Subtracted No.	Energy (keV)	Original Area	Orig. Uncert.	Area Background	Ambient Uncert.	Backgr. Uncert.	Subtracted Area
1	238.55	7.17E+001	17.15				7.17E+001
1.72E+001							
2	582.92	1.50E+001	8.71				1.50E+001
8.71E+000							
3	661.33	3.15E+001	8.75				3.15E+001
8.75E+000							
4	1172.72	1.89E+002	16.16				1.89E+002
1.62E+001							
5	1332.20	1.38E+002	12.43				1.38E+002
1.24E+001							
6	1460.20	3.25E+001	6.03				3.25E+001
6.03E+000							

M = First peak in a multiplet region
m = Other peak in a multiplet region
F = Fitted singlet

Errors quoted at 1.000 sigma

**

 N U C L I D E M D A R E P O R T

**

Detector Name: DET01
Sample Geometry: 1-L Marinelli
Sample Title: 1-L Solid Releas
Nuclide Library Used: C:\GENIE2K\CAMFILES\AnSolMDA.NLB

Activity (uCi/gram)	Nuclide Name	Energy (keV)	Yield (%)	Line MDA (uCi/gram)	Nuclide MDA (uCi/gram)	
4.73E-007	CS-134	563.23	8.38	1.14E-006	1.14E-007	-
7.60E-008		569.32	15.43	8.02E-007		
5.15E-008		604.70	97.60	1.14E-007		-
8.11E-009		795.84	85.40	2.16E-007		
2.08E-008		801.93	8.73	1.96E-006		-
7.51E-007		1365.15	3.04	3.53E-006		
1.46E-007	+ CS-137	661.65*	85.12	1.57E-007	1.57E-007	

+ = Nuclide identified during the nuclide identification
* = Energy line found in the spectrum
> = MDA value not calculated
@ = Half-life too short to be able to perform the decay
correction

*
***** N U C L I D E I D E N T I F I C A T I O N R E P O R T

*

Sample Title: 1-L Solid Releas
Nuclide Library Used: C:\GENIE2K\CAMFILES\AnSolMDA.NLB

..... IDENTIFIED NUCLIDES

Nuclide Name	Id Confidence	Energy (keV)	Yield (%)	Activity (uCi/gram)	Activity Uncertainty
K-40	0.942	1460.81*	10.67	2.46E-006	4.60E-007
CO-60	0.975	1173.22*	100.00	1.25E-006	1.10E-007
		1332.49*	100.00	1.03E-006	9.48E-008
CS-137	0.984	661.65*	85.12	1.46E-007	4.07E-008
PB-212	0.709	87.20	6.30		
		89.80	1.75		
		238.63*	44.60	2.30E-007	5.52E-008
		300.09	3.41		

* = Energy line found in the spectrum.
@ = Energy line not used for Weighted Mean Activity
Energy Tolerance : 1.000 keV
Nuclide confidence index threshold = 0.30
Errors quoted at 1.000 sigma

 *
 ***** INTERFERENCE CORRECTED REPORT *****

 *

Nuclide Name	Nuclide Id Confidence	Wt mean Activity (uCi/gram)	Wt mean Activity Uncertainty
K-40	0.942	2.46E-006	4.60E-007
CO-60	0.975	1.13E-006	7.18E-008
CS-137	0.984	1.46E-007	4.07E-008
PB-212	0.709	2.30E-007	5.52E-008
Total Activity		3.96E-006	

? = nuclide is part of an undetermined solution
 X = nuclide rejected by the interference analysis
 @ = nuclide contains energy lines not used in Weighted Mean

Activity

Errors quoted at 1.000 sigma

***** UNIDENTIFIED PEAKS *****

Peak Locate Performed on: 4/18/09 1:36:19 PM
 Peak Locate From Channel: 100
 Peak Locate To Channel: 4096

Peak No.	Energy (keV)	Peak Size in Counts per Second	Peak CPS % Uncertainty
2	582.92	8.35E-003	58.07

M = First peak in a multiplet region
 m = Other peak in a multiplet region
 F = Fitted singlet

Errors quoted at 1.000 sigma