

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042924\
 Data File : BG061053.D
 Acq On : 29 Apr 2024 19:06
 Operator : MA/JU
 Sample : P2230-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SOIL

Quant Time: Apr 30 01:22:08 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 22 04:22:00 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.941	152	22332	20.000	ng	0.00
21) Naphthalene-d8	10.737	136	86939	20.000	ng	# 0.00
39) Acenaphthene-d10	14.568	164	54220	20.000	ng	0.00
64) Phenanthrene-d10	17.318	188	199760	20.000	ng	# 0.00
76) Chrysene-d12	21.572	240	185329	20.000	ng	0.00
86) Perylene-d12	24.697	264	24873	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.532	112	145522	131.808	ng	0.00
7) Phenol-d6	7.112	99	159851	97.583	ng	0.00
23) Nitrobenzene-d5	9.098	82	157248	81.264	ng	0.00
42) 2,4,6-Tribromophenol	16.061	330	215489	180.987	ng	0.00
45) 2-Fluorobiphenyl	13.193	172	463395	118.285	ng	0.00
79) Terphenyl-d14	19.938	244	1122015	93.975	ng	0.00
Target Compounds						
59) 2,3,4,6-Tetrachlorophenol	15.197	232	29812	19.715	ng	97
70) Pentachlorophenol	16.977	266	787235	412.274	ng	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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