

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050924\
 Data File : BG061155.D
 Acq On : 9 May 2024 17:32
 Operator : MA/JU
 Sample : P2390-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PE-4

Quant Time: May 09 18:15:00 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG043024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 05:11:25 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.935	152	25611	20.000	ng	-0.03
21) Naphthalene-d8	10.732	136	96022	20.000	ng	#-0.03
39) Acenaphthene-d10	14.563	164	71419	20.000	ng	-0.03
64) Phenanthrene-d10	17.312	188	158877	20.000	ng	#-0.02
76) Chrysene-d12	21.572	240	168595	20.000	ng	-0.02
86) Perylene-d12	24.692	264	206918	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.526	112	104914	83.311	ng	-0.02
7) Phenol-d6	7.107	99	141187	75.788	ng	-0.03
23) Nitrobenzene-d5	9.087	82	106316	54.887	ng	-0.03
42) 2,4,6-Tribromophenol	16.061	330	102062	71.495	ng	-0.01
45) 2-Fluorobiphenyl	13.182	172	250763	51.920	ng	-0.03
79) Terphenyl-d14	19.933	244	457883	45.169	ng	-0.02
Target Compounds						
						Qvalue
75) Fluoranthene	19.369	202	23485	2.360	ng	91
78) Pyrene	19.733	202	25158	2.347	ng	# 91

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

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