

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110722\
 Data File : BG055492.D
 Acq On : 8 Nov 2022 1:01
 Operator : CG/JU
 Sample : N5415-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 R-22

Quant Time: Nov 08 03:55:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 03:46:08 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.078	152	28156	20.000	ng	0.00
21) Naphthalene-d8	10.910	136	129961	20.000	ng	# 0.00
39) Acenaphthene-d10	14.717	164	91605	20.000	ng	-0.01
64) Phenanthrene-d10	17.455	188	198541	20.000	ng	-0.01
76) Chrysene-d12	21.715	240	192894	20.000	ng	-0.01
86) Perylene-d12	24.970	264	217256	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.605	112	162440	103.608	ng	0.00
7) Phenol-d6	7.244	99	220209	101.283	ng	0.00
23) Nitrobenzene-d5	9.259	82	124415	51.021	ng	0.00
42) 2,4,6-Tribromophenol	16.204	330	127386	101.948	ng	-0.01
45) 2-Fluorobiphenyl	13.343	172	355661	57.565	ng	0.00
79) Terphenyl-d14	20.047	244	578340	58.302	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.500	202	31272	2.328	ng	97
78) Pyrene	19.859	202	31028	2.280	ng	99
88) Benzo(b)fluoranthene	23.930	252	28354	2.055	ng	# 90

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

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