

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102122\
 Data File : BG055246.D
 Acq On : 21 Oct 2022 23:15
 Operator : CG/JU
 Sample : N5217-16
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-K1-COMP

Quant Time: Oct 22 01:33:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG101922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 20 09:21:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.287	152	52913	20.000	ng	0.00
21) Naphthalene-d8	11.113	136	260013	20.000	ng	0.00
39) Acenaphthene-d10	14.897	164	176329	20.000	ng	0.00
64) Phenanthrene-d10	17.629	188	343170	20.000	ng	0.00
76) Chrysene-d12	21.907	240	226858	20.000	ng	-0.01
86) Perylene-d12	25.309	264	237508	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.808	112	411784	140.014	ng	0.00
7) Phenol-d6	7.424	99	580775	129.595	ng	0.00
23) Nitrobenzene-d5	9.457	82	358476	66.560	ng	0.00
42) 2,4,6-Tribromophenol	16.372	330	205265	112.343	ng	0.00
45) 2-Fluorobiphenyl	13.522	172	767669	69.541	ng	0.00
79) Terphenyl-d14	20.203	244	829046	68.094	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.656	202	63354	3.123	ng	96
78) Pyrene	20.021	202	63780	3.731	ng	98
81) Benzo(a)anthracene	21.883	228	29488	2.004	ng	95
83) Chrysene	21.954	228	27589	2.013	ng	96
88) Benzo(b)fluoranthene	24.222	252	41601	2.950	ng	# 90
90) Benzo(a)pyrene	25.144	252	31471	2.600	ng	# 90

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

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