

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031022\
 Data File : BG052631.D
 Acq On : 10 Mar 2022 13:07
 Operator : CG/JU
 Sample : N1845-13RE 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SAMPLE-13RE

Quant Time: Mar 11 03:44:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG030922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 10 06:37:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.170	152	43598	20.000	ng	0.00
21) Naphthalene-d8	11.007	136	170753	20.000	ng	0.00
39) Acenaphthene-d10	14.826	164	109570	20.000	ng	0.00
64) Phenanthrene-d10	17.575	188	227915	20.000	ng	0.00
76) Chrysene-d12	21.881	240	224386	20.000	ng	0.00
86) Perylene-d12	25.306	264	232754	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.703	112	17214	6.821	ng	0.00
7) Phenol-d6	7.318	99	79269	21.702	ng	0.00
23) Nitrobenzene-d5	9.345	82	76093	20.201	ng	0.00
42) 2,4,6-Tribromophenol	16.318	330	1695	1.091	ng	0.00
45) 2-Fluorobiphenyl	13.445	172	166084	20.459	ng	0.00
79) Terphenyl-d14	20.165	244	252948	19.948	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.616	178	26539	2.219	ng	96
75) Fluoranthene	19.619	202	47021	3.149	ng	100
78) Pyrene	19.983	202	42806	2.854	ng	98

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

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