

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042322\
 Data File : BG053291.D
 Acq On : 23 Apr 2022 13:48
 Operator : CG/JU
 Sample : N2526-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MH-3S

Quant Time: Apr 24 02:49:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Apr 24 01:13:30 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.105	152	30192	20.000	ng	0.00
21) Naphthalene-d8	10.919	136	126953	20.000	ng	0.00
39) Acenaphthene-d10	14.732	164	100352	20.000	ng	0.00
64) Phenanthrene-d10	17.475	188	239340	20.000	ng	0.00
76) Chrysene-d12	21.757	240	230147	20.000	ng	0.00
86) Perylene-d12	25.053	264	236142	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.673	112	185191	105.144	ng	0.00
7) Phenol-d6	7.265	99	269308	99.152	ng	0.00
23) Nitrobenzene-d5	9.268	82	192850	61.679	ng	0.00
42) 2,4,6-Tribromophenol	16.218	330	144228	90.337	ng	0.00
45) 2-Fluorobiphenyl	13.351	172	450029	61.837	ng	0.00
79) Terphenyl-d14	20.077	244	841329	61.740	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.516	178	28399	2.172	ng	94
75) Fluoranthene	19.525	202	42366	2.528	ng	93
78) Pyrene	19.889	202	49053	2.993	ng	98
88) Benzo(b)fluoranthene	24.001	252	39634	2.660	ng #	93
90) Benzo(a)pyrene	24.894	252	30674	2.417	ng #	92

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

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