

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042624\
 Data File : BF137735.D
 Acq On : 26 Apr 2024 15:16
 Operator : CG\JU
 Sample : P2218-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-131-GW-01

Quant Time: Apr 26 15:38:22 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 24 03:57:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.057	152	263842	20.000	ng	0.00
21) Naphthalene-d8	8.345	136	1035157	20.000	ng	0.00
39) Acenaphthene-d10	10.104	164	578632	20.000	ng	-0.01
64) Phenanthrene-d10	11.598	188	973877	20.000	ng	0.00
76) Chrysene-d12	14.245	240	496435	20.000	ng	-0.01
86) Perylene-d12	15.821	264	495411	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.675	112	793107	45.662	ng	0.00
7) Phenol-d6	6.663	99	648813	29.416	ng	-0.01
23) Nitrobenzene-d5	7.622	82	1251771	64.191	ng	0.00
42) 2,4,6-Tribromophenol	10.898	330	534657	91.569	ng	0.00
45) 2-Fluorobiphenyl	9.416	172	2430060	62.752	ng	-0.01
79) Terphenyl-d14	13.186	244	2455536	79.614	ng	-0.01
Target Compounds						
						Qvalue
38) 1-Methylnaphthalene	9.157	142	91219	2.691	ng	99
52) Acenaphthene	10.139	154	232503	7.096	ng	99
55) Dibenzofuran	10.310	168	190902	3.975	ng	98
58) Fluorene	10.651	166	185897	4.953	ng	98
71) Phenanthrene	11.622	178	158574	3.076	ng	97

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

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