

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042624\
 Data File : BF137736.D
 Acq On : 26 Apr 2024 15:47
 Operator : CG\JU
 Sample : P2218-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-131-GW-02

Quant Time: Apr 26 16:22:22 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	257405	20.000	ng	0.00
21) Naphthalene-d8	8.345	136	1024827	20.000	ng	0.00
39) Acenaphthene-d10	10.104	164	571802	20.000	ng	-0.01
64) Phenanthrene-d10	11.598	188	976845	20.000	ng	0.00
76) Chrysene-d12	14.245	240	514641	20.000	ng	-0.01
86) Perylene-d12	15.821	264	512335	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.675	112	808323	47.702	ng	0.00
7) Phenol-d6	6.663	99	661592	30.745	ng	-0.01
23) Nitrobenzene-d5	7.622	82	1225224	63.463	ng	0.00
42) 2,4,6-Tribromophenol	10.898	330	557931	96.697	ng	0.00
45) 2-Fluorobiphenyl	9.416	172	2379773	62.187	ng	-0.01
79) Terphenyl-d14	13.186	244	2522448	78.891	ng	-0.01
Target Compounds						
						Qvalue
38) 1-Methylnaphthalene	9.157	142	115246	3.434	ng	98
52) Acenaphthene	10.139	154	293215	9.056	ng	98
55) Dibenzofuran	10.310	168	241617	5.091	ng	97
58) Fluorene	10.651	166	244004	6.579	ng	99
71) Phenanthrene	11.622	178	208017	4.023	ng	97

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

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