

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120823\
 Data File : BF136469.D
 Acq On : 08 Dec 2023 16:04
 Operator : CG\JU
 Sample : 05629-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ-216

Quant Time: Dec 08 16:27:03 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 06 09:40:27 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.792	152	119423	20.000	ng	-0.01
21) Naphthalene-d8	8.069	136	485790	20.000	ng	-0.01
39) Acenaphthene-d10	9.827	164	249652	20.000	ng	-0.01
64) Phenanthrene-d10	11.321	188	463735	20.000	ng	0.00
76) Chrysene-d12	13.968	240	305088	20.000	ng	#-0.01
86) Perylene-d12	15.451	264	250443	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.451	112	939805	116.220	ng	0.02
7) Phenol-d6	6.445	99	1091890	108.225	ng	0.00
23) Nitrobenzene-d5	7.357	82	791920	88.128	ng	-0.01
42) 2,4,6-Tribromophenol	10.621	330	383715	124.654	ng	-0.01
45) 2-Fluorobiphenyl	9.151	172	1569316	85.773	ng	-0.01
79) Terphenyl-d14	12.921	244	1688091	72.929	ng	0.00
Target Compounds						
31) Naphthalene	8.092	128	354808	13.244	ng	98
37) 2-Methylnaphthalene	8.780	142	37568	2.204	ng	87
52) Acenaphthene	9.857	154	42096	2.789	ng	100
55) Dibenzofuran	10.033	168	45723	2.074	ng	95
58) Fluorene	10.374	166	38528	2.177	ng	92
71) Phenanthrene	11.339	178	60027	2.355	ng	99
73) Carbazole	11.551	167	63987	2.816	ng	98
77) Benzidine	12.663	184	77934	17.177	ng	99

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

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