

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062521\
 Data File : BF124454.D
 Acq On : 25 Jun 2021 16:27
 Operator : JU/CG
 Sample : M2812-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 293-347

Quant Time: Jun 25 16:55:17 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF062321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 25 13:23:23 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.704	152	32950	20.000	ng	0.00
21) Naphthalene-d8	7.986	136	121126	20.000	ng	# 0.00
39) Acenaphthene-d10	9.733	164	69684	20.000	ng	0.00
64) Phenanthrene-d10	11.222	188	111726	20.000	ng	0.00
76) Chrysene-d12	13.863	240	80581	20.000	ng	0.00
86) Perylene-d12	15.286	264	80401	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	264802	126.389	ng	0.01
7) Phenol-d6	6.363	99	308981	112.948	ng	0.00
23) Nitrobenzene-d5	7.275	82	252441	86.413	ng	0.00
42) 2,4,6-Tribromophenol	10.533	330	116050	139.393	ng	0.00
45) 2-Fluorobiphenyl	9.063	172	497924	87.531	ng	0.00
79) Terphenyl-d14	12.810	244	443764	82.484	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.010	128	693549	97.323	ng	98
37) 2-Methylnaphthalene	8.698	142	70013	13.884	ng	96
38) 1-Methylnaphthalene	8.792	142	43825	9.343	ng	96
46) 1,1'-Biphenyl	9.163	154	15217	2.569	ng	94
52) Acenaphthene	9.769	154	30706	6.906	ng	90
55) Dibenzofuran	9.945	168	31763	4.482	ng	90
58) Fluorene	10.286	166	26549	4.875	ng	86
71) Phenanthrene	11.245	178	33302	4.728	ng	98
73) Carbazole	11.457	167	28760	4.702	ng	# 97

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

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