

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062521\
 Data File : BF124453.D
 Acq On : 25 Jun 2021 15:55
 Operator : JU/CG
 Sample : M2812-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 356-349

Quant Time: Jun 25 16:54:57 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	32610	20.000	ng	0.00
21) Naphthalene-d8	7.987	136	123015	20.000	ng	# 0.00
39) Acenaphthene-d10	9.739	164	69776	20.000	ng	0.00
64) Phenanthrene-d10	11.222	188	119867	20.000	ng	0.00
76) Chrysene-d12	13.863	240	85400	20.000	ng	0.00
86) Perylene-d12	15.292	264	91324	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	270182	130.301	ng	0.01
7) Phenol-d6	6.363	99	304582	112.501	ng	0.00
23) Nitrobenzene-d5	7.275	82	260566	87.825	ng	0.00
42) 2,4,6-Tribromophenol	10.533	330	116731	140.026	ng	0.00
45) 2-Fluorobiphenyl	9.063	172	513928	90.225	ng	0.00
79) Terphenyl-d14	12.810	244	468084	82.095	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.375	94	5907	2.018	ng	77
31) Naphthalene	8.016	128	1274294	176.071	ng	98
37) 2-Methylnaphthalene	8.698	142	91759	17.916	ng	87
38) 1-Methylnaphthalene	8.798	142	53089	11.145	ng	97
46) 1,1'-Biphenyl	9.163	154	20827	3.512	ng	94
52) Acenaphthene	9.769	154	50660	11.378	ng	95
55) Dibenzofuran	9.945	168	54906	7.737	ng	96
58) Fluorene	10.286	166	41668	7.642	ng	99
71) Phenanthrene	11.245	178	53801	7.120	ng	97
73) Carbazole	11.463	167	66942	10.202	ng	98

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

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