

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030221\
 Data File : BF123337.D
 Acq On : 2 Mar 2021 15:39
 Operator : JU/CG
 Sample : M1500-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T-2

Quant Time: Mar 02 17:18:26 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF021921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 19 16:08:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.851	152	135637	20.00	ng	# 0.00
21) Naphthalene-d8	8.145	136	445475	20.00	ng	# 0.00
39) Acenaphthene-d10	9.910	164	277840	20.00	ng	0.00
64) Phenanthrene-d10	11.398	188	488550	20.00	ng	# 0.00
76) Chrysene-d12	14.039	240	344619	20.00	ng	0.00
86) Perylene-d12	15.515	264	300130	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	805645	85.50	ng	0.00
7) Phenol-d6	6.487	99	1048011	81.86	ng	0.00
23) Nitrobenzene-d5	7.416	82	711759	73.14	ng	-0.01
42) 2,4,6-Tribromophenol	10.704	330	245803	86.88	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	1076555	55.66	ng	0.00
79) Terphenyl-d14	12.986	244	1091510	60.90	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.498	94	29906	2.14	ng	86
22) Acetophenone	7.251	105	350006	30.29	ng	# 56
37) 2-Methylnaphthalene	8.857	142	114265	7.03	ng	99
38) 1-Methylnaphthalene	8.957	142	434580	28.57	ng	99
46) 1,1'-Biphenyl	9.322	154	91610	3.97	ng	94
52) Acenaphthene	9.939	154	85847	4.51	ng	# 83
58) Fluorene	10.457	166	176759	8.72	ng	# 90
71) Phenanthrene	11.427	178	555151	19.22	ng	96
78) Pyrene	12.845	202	128418	4.77	ng	# 95

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

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