

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111524\
 Data File : BF140395.D
 Acq On : 15 Nov 2024 11:33
 Operator : RC/JU
 Sample : P4799-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-TA2-02-C

Quant Time: Nov 15 12:08:27 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF111324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 14:40:06 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	126128	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	478686	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	258727	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	489102	20.000	ng	0.00
76) Chrysene-d12	14.062	240	278626	20.000	ng	0.01
86) Perylene-d12	15.586	264	287340	20.000	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	939985	127.245	ng	0.03
7) Phenol-d6	6.522	99	1219220	121.819	ng	0.02
23) Nitrobenzene-d5	7.434	82	784930	85.436	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	363116	141.134	ng	0.00
45) 2-Fluorobiphenyl	9.227	172	1364655	84.790	ng	0.00
79) Terphenyl-d14	12.992	244	1584307	98.700	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.175	128	770176	31.717	ng	99
37) 2-Methylnaphthalene	8.869	142	103203	6.628	ng	99
38) 1-Methylnaphthalene	8.963	142	164098	10.762	ng	99
52) Acenaphthene	9.945	154	184909	12.620	ng	99
58) Fluorene	10.457	166	94791	5.785	ng	99
71) Phenanthrene	11.421	178	208137	9.059	ng	99
77) Benizidine	12.739	184	80699	7.546	ng	99

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

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