

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111524\
 Data File : BF140398.D
 Acq On : 15 Nov 2024 12:52
 Operator : RC/JU
 Sample : P4799-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-TA2-03-C

Quant Time: Nov 15 13:18:49 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	117399	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	435226	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	234558	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	431505	20.000	ng	0.00
76) Chrysene-d12	14.051	240	249173	20.000	ng	0.00
86) Perylene-d12	15.557	264	204971	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	879744	127.945	ng	0.03
7) Phenol-d6	6.522	99	1141001	122.480	ng	0.02
23) Nitrobenzene-d5	7.434	82	739645	88.546	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	326513	139.984	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	1304219	89.385	ng	0.00
79) Terphenyl-d14	12.986	244	1458423	101.597	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.175	128	133494	6.046	ng	99
37) 2-Methylnaphthalene	8.869	142	28721	2.029	ng	95
38) 1-Methylnaphthalene	8.963	142	45614	3.290	ng	99
52) Acenaphthene	9.939	154	49484	3.725	ng	98
58) Fluorene	10.457	166	30129	2.028	ng	97
71) Phenanthrene	11.422	178	81919	4.041	ng	98

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

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