

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111524\
 Data File : BF140399.D
 Acq On : 15 Nov 2024 13:18
 Operator : RC/JU
 Sample : P4799-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-TA1-01-C

Quant Time: Nov 15 14:04:17 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	127141	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	473766	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	259206	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	478884	20.000	ng	0.00
76) Chrysene-d12	14.051	240	265821	20.000	ng	0.00
86) Perylene-d12	15.563	264	271820	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	871858	117.082	ng	0.03
7) Phenol-d6	6.516	99	1093135	108.351	ng	0.01
23) Nitrobenzene-d5	7.434	82	788403	86.705	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	344401	133.613	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	1409740	87.429	ng	0.00
79) Terphenyl-d14	12.986	244	1534702	100.215	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.181	128	2114875	87.998	ng	100
37) 2-Methylnaphthalene	8.863	142	286408	18.585	ng	99
38) 1-Methylnaphthalene	8.963	142	338220	22.412	ng	100
52) Acenaphthene	9.939	154	87029	5.929	ng	98
58) Fluorene	10.457	166	49290	3.003	ng	98
71) Phenanthrene	11.422	178	90172	4.008	ng	100

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

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