

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120624\
 Data File : BF140763.D
 Acq On : 06 Dec 2024 13:48
 Operator : RC/JU
 Sample : P5133-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MOO-24-00374

Quant Time: Dec 06 23:26:00 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 21 15:23:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	58120	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	154680	20.000	ng	# 0.00
39) Acenaphthene-d10	9.928	164	96966	20.000	ng	0.01
64) Phenanthrene-d10	11.422	188	149935	20.000	ng	# 0.02
76) Chrysene-d12	14.086	240	120887	20.000	ng	# 0.04
86) Perylene-d12	15.604	264	110610	20.000	ng	0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	169241	49.684	ng	0.00
7) Phenol-d6	6.516	99	205729	45.684	ng	0.00
23) Nitrobenzene-d5	7.434	82	134918	44.615	ng	0.00
42) 2,4,6-Tribromophenol	10.722	330	38693	37.310	ng	0.02
45) 2-Fluorobiphenyl	9.239	172	201545	30.969	ng	0.00
79) Terphenyl-d14	13.016	244	214084	27.576	ng	0.02
Target Compounds						Qvalue
31) Naphthalene	8.181	128	196863	24.709	ng	# 69
37) 2-Methylnaphthalene	8.881	142	331297	65.467	ng	99
38) 1-Methylnaphthalene	8.981	142	198463	40.010	ng	92
46) 1,1'-Biphenyl	9.339	154	103128	14.245	ng	96
71) Phenanthrene	11.445	178	57524	7.981	ng	# 61
78) Pyrene	12.892	202	52959	4.738	ng	# 78
84) Bis(2-ethylhexyl)phtha...	14.045	149	83246	16.373	ng	# 73

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

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