

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120121\
 Data File : BP008181.D
 Acq On : 01 Dec 2021 19:52
 Operator : CG/JU
 Sample : M4721-01 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SAMPLE-1

Quant Time: Dec 01 21:16:30 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 01 12:06:05 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	87857	20.000	ng	0.00
21) Naphthalene-d8	10.593	136	312105	20.000	ng	0.00
39) Acenaphthene-d10	14.440	164	201288	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	454417	20.000	ng	0.00
76) Chrysene-d12	21.304	240	514135	20.000	ng	0.00
86) Perylene-d12	23.698	264	565286	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	198023	36.290	ng	0.00
7) Phenol-d6	6.981	99	243529	33.061	ng	0.00
23) Nitrobenzene-d5	8.958	82	166511	24.260	ng	0.00
42) 2,4,6-Tribromophenol	15.940	330	100949	39.237	ng	0.00
45) 2-Fluorobiphenyl	13.063	172	326601	23.568	ng	0.00
79) Terphenyl-d14	19.775	244	575774	23.405	ng	0.00

Target Compounds	Qvalue

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

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