

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121021\
 Data File : BP008349.D
 Acq On : 10 Dec 2021 19:59
 Operator : CG/JU
 Sample : M4966-23
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P3-COMP

Quant Time: Dec 11 04:01:11 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 11 03:53:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.758	152	74839	20.000	ng	0.00
21) Naphthalene-d8	10.557	136	269479	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	163236	20.000	ng	0.00
64) Phenanthrene-d10	17.181	188	318353	20.000	ng	0.00
76) Chrysene-d12	21.286	240	262878	20.000	ng	0.00
86) Perylene-d12	23.669	264	296330	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	582612	125.344	ng	0.00
7) Phenol-d6	6.952	99	665720	106.097	ng	0.00
23) Nitrobenzene-d5	8.922	82	553273	93.362	ng	0.00
42) 2,4,6-Tribromophenol	15.916	330	276373	132.462	ng	0.00
45) 2-Fluorobiphenyl	13.028	172	1035203	92.116	ng	0.00
79) Terphenyl-d14	19.751	244	1343142	106.782	ng	0.00
Target Compounds						
84) Bis(2-ethylhexyl)phtha...	21.198	149	58496	5.147	ng	Qvalue 98

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

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