

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120723\
 Data File : BP018683.D
 Acq On : 07 Dec 2023 20:51
 Operator : MA/JU
 Sample : 05459-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA2

Quant Time: Dec 07 21:52:29 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 04 21:39:40 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	194526	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.640	136	717149	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.475	164	452006	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.222	188	1109699	20.000	ng/u1	0.00
79) Chrysene-d12	21.310	240	1359469	20.000	ng/u1	0.00
88) Perylene-d12	23.680	264	1899524	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	5314	0.931	ng/uL	0.00
4) Pyridine-d5	3.687	84	56856	3.718	ng/u1	0.00
7) Phenol-d5	7.010	99	49958	2.559	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.181	67	34143	2.951	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.375	132	40881	2.684	ng/u1	-0.01
15) 4-Methylphenol-d8	8.557	113	17515	1.109	ng/u1	-0.01
21) Nitrobenzene-d5	9.004	128	18729	2.934	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.728	143	17641	2.666	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.263	165	34857	2.576	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.781	131	25331	1.432	ng/u1	-0.01
46) Dimethylphthalate-d6	13.892	166	125576	3.113	ng/u1	-0.01
49) Acenaphthylene-d8	14.169	160	131514	2.961	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.681	143	9298	1.404	ng/u1	0.00
60) Fluorene-d10	15.469	176	115427	3.409	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.586	200	7056	1.060	ng/u1	0.00
73) Anthracene-d10	17.322	188	175633	3.011	ng/u1	0.00
81) Pyrene-d10	19.557	212	291055	2.750	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.521	264	349678	3.157	ng/u1	-0.01
Target Compounds						
16) Acetophenone	8.669	105	89477	3.800	ng/u1	96

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

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