

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120723\
 Data File : BP018677.D
 Acq On : 07 Dec 2023 17:27
 Operator : MA/JU
 Sample : 05459-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AJ0

Quant Time: Dec 07 21:39:39 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	195417	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.640	136	739616	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.475	164	477563	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.222	188	1168544	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	1369550	20.000	ng/ul	0.00
88) Perylene-d12	23.680	264	1688912	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	19290	3.364	ng/uL	0.00
4) Pyridine-d5	3.681	84	250319	16.296	ng/ul	0.00
7) Phenol-d5	7.010	99	310099	15.813	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.181	67	182183	15.674	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.375	132	240993	15.749	ng/ul	-0.01
15) 4-Methylphenol-d8	8.557	113	240192	15.134	ng/ul	-0.01
21) Nitrobenzene-d5	9.004	128	110541	16.792	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.728	143	117503	17.217	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.263	165	233122	16.705	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.781	131	313302	17.178	ng/ul	-0.01
46) Dimethylphthalate-d6	13.892	166	736728	17.285	ng/ul	-0.01
49) Acenaphthylene-d8	14.169	160	809732	17.253	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.675	143	35536	5.079	ng/ul	-0.01
60) Fluorene-d10	15.469	176	655905	18.333	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	25854	3.690	ng/ul	-0.01
73) Anthracene-d10	17.322	188	1065212	17.340	ng/ul	0.00
81) Pyrene-d10	19.557	212	1446754	13.571	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.527	264	1707528	17.340	ng/ul	0.00

Target Compounds Qvalue

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

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