

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017334.D
 Acq On : 25 Sep 2023 22:42
 Operator : MA/JU
 Sample : 04429-19
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHE3

Quant Time: Sep 26 00:22:05 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 21 23:51:07 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	215835	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.751	136	918459	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.586	164	576209	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.357	188	1273085	20.000	ng/ul	-0.01
79) Chrysene-d12	21.463	240	1177932	20.000	ng/ul	-0.01
88) Perylene-d12	23.974	264	1307657	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	24087	4.584	ng/uL	0.00
4) Pyridine-d5	3.734	84	141192	9.611	ng/ul	0.00
7) Phenol-d5	7.087	99	117199	6.612	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.269	67	374305	34.989	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.452	132	328619	23.374	ng/ul	-0.01
15) 4-Methylphenol-d8	8.640	113	215341	15.625	ng/ul	-0.02
21) Nitrobenzene-d5	9.128	128	240575	36.232	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.840	143	211789	29.375	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.363	165	369762	27.146	ng/ul	-0.02
31) 4-Chloroaniline-d4	10.916	131	417469	21.158	ng/ul	-0.01
46) Dimethylphthalate-d6	14.004	166	1610517	39.016	ng/ul	-0.01
49) Acenaphthylene-d8	14.286	160	1722240	36.277	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.828	143	34688	4.824	ng/ul	0.00
60) Fluorene-d10	15.586	176	1391981	39.171	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.716	200	188002	26.201	ng/ul	-0.01
73) Anthracene-d10	17.457	188	2469677	43.740	ng/ul	-0.01
81) Pyrene-d10	19.698	212	3046338	47.824	ng/ul	-0.02
92) Benzo(a)pyrene-d12	23.810	264	2941109	46.714	ng/ul	-0.02

Target Compounds Qvalue

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

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