

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092923\
 Data File : BP017456.D
 Acq On : 29 Sep 2023 21:17
 Operator : MA/JU
 Sample : 04497-18
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHM6

Quant Time: Sep 29 22:37:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 29 22:31:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	224146	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.787	136	932607	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.646	164	592032	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.434	188	1341915	20.000	ng/ul	-0.01
79) Chrysene-d12	21.563	240	1310143	20.000	ng/ul	-0.01
88) Perylene-d12	24.133	264	1460233	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	8956	1.641	ng/uL	0.00
4) Pyridine-d5	3.734	84	118145	7.744	ng/ul	0.00
7) Phenol-d5	7.105	99	126156	6.853	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.293	67	188321	16.951	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.469	132	199443	13.660	ng/ul	-0.01
15) 4-Methylphenol-d8	8.669	113	165404	11.557	ng/ul	-0.02
21) Nitrobenzene-d5	9.158	128	130768	19.395	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.875	143	117361	16.031	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.405	165	216825	15.677	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.958	131	315882	15.766	ng/ul	-0.01
46) Dimethylphthalate-d6	14.057	166	812162	19.150	ng/ul	-0.01
49) Acenaphthylene-d8	14.340	160	941830	19.308	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.887	143	17621	2.385	ng/ul	0.00
60) Fluorene-d10	15.645	176	726200	19.889	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.793	200	16427	2.172	ng/ul	-0.01
73) Anthracene-d10	17.534	188	1196313	20.101	ng/ul	-0.01
81) Pyrene-d10	19.781	212	1431912	20.211	ng/ul	-0.02
92) Benzo(a)pyrene-d12	23.963	264	1382630	19.666	ng/ul	-0.02

Target Compounds Qvalue

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

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