

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090723\
 Data File : BP017003.D
 Acq On : 07 Sep 2023 23:55
 Operator : MA/JU
 Sample : 04166-18
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJY0

Quant Time: Sep 08 00:42:02 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 06 01:52:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	400564	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	1755520	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	1074399	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	2251693	20.000	ng/u1	0.00
79) Chrysene-d12	21.498	240	1679777	20.000	ng/u1	-0.01
88) Perylene-d12	24.033	264	1594760	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	51695	4.882	ng/uL	0.00
4) Pyridine-d5	3.764	84	326660	10.796	ng/u1	0.00
7) Phenol-d5	7.122	99	326079	8.910	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	933911	40.799	ng/u1	0.00
11) 2-Chlorophenol-d4	7.499	132	846041	31.943	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	591750	21.225	ng/u1	-0.01
21) Nitrobenzene-d5	9.175	128	543845	39.931	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	575645	38.991	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.416	165	954730	35.028	ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	1319812	33.562	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	3474728	43.447	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	3786294	41.023	ng/u1	0.00
54) 4-Nitrophenol-d4	14.845	143	113155	7.243	ng/u1	-0.01
60) Fluorene-d10	15.628	176	2914046	42.557	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	471131	34.768	ng/u1	0.00
73) Anthracene-d10	17.504	188	4580464	44.295	ng/u1	0.00
81) Pyrene-d10	19.739	212	4954225	48.421	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.868	264	3732107	46.129	ng/u1	-0.01

Target Compounds Qvalue

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

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