

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017333.D
 Acq On : 25 Sep 2023 22:08
 Operator : MA/JU
 Sample : 04429-18
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJ52

Quant Time: Sep 26 00:41:22 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	176984	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.752	136	772823	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.587	164	474335	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.357	188	1041464	20.000	ng/u1	-0.01
79) Chrysene-d12	21.463	240	931166	20.000	ng/u1	-0.01
88) Perylene-d12	23.980	264	1028117	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	22382	5.194	ng/uL	0.00
4) Pyridine-d5	3.734	84	192763	16.002	ng/u1	0.00
7) Phenol-d5	7.087	99	119121	8.195	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.269	67	374624	42.706	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.446	132	363844	31.560	ng/u1	-0.02
15) 4-Methylphenol-d8	8.640	113	229900	20.344	ng/u1	-0.02
21) Nitrobenzene-d5	9.128	128	241540	43.232	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.840	143	250519	41.295	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.363	165	436309	38.068	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.916	131	647782	39.017	ng/u1	-0.01
46) Dimethylphthalate-d6	14.004	166	1649015	48.529	ng/u1	-0.01
49) Acenaphthylene-d8	14.287	160	1761237	45.066	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.822	143	32712	5.526	ng/u1	0.00
60) Fluorene-d10	15.587	176	1421604	48.596	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.716	200	198515	33.819	ng/u1	-0.01
73) Anthracene-d10	17.457	188	2324283	50.320	ng/u1	-0.01
81) Pyrene-d10	19.704	212	2733399	54.283	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.816	264	2588643	52.295	ng/u1	-0.02

Target Compounds Qvalue

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

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