

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
 Data File : BP017305.D
 Acq On : 23 Sep 2023 01:43
 Operator : MA/JU
 Sample : 04410-20
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHH9

Quant Time: Sep 23 02:17:13 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	148370	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.752	136	602836	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.593	164	362760	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.363	188	815120	20.000	ng/u1	0.00
79) Chrysene-d12	21.469	240	779887	20.000	ng/u1	0.00
88) Perylene-d12	23.980	264	859625	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	19596	5.425	ng/uL	0.00
4) Pyridine-d5	3.746	84	86403	8.556	ng/u1	0.00
7) Phenol-d5	7.087	99	91083	7.475	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.275	67	233918	31.809	ng/u1	0.00
11) 2-Chlorophenol-d4	7.452	132	249445	25.810	ng/u1	-0.01
15) 4-Methylphenol-d8	8.646	113	155539	16.418	ng/u1	-0.01
21) Nitrobenzene-d5	9.128	128	145759	33.445	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.846	143	153642	32.468	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	275062	30.766	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.922	131	102206	7.892	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	1002460	38.575	ng/u1	-0.01
49) Acenaphthylene-d8	14.293	160	1026105	34.331	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	27738	6.127	ng/u1	0.00
60) Fluorene-d10	15.587	176	856295	38.275	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.722	200	121795	26.511	ng/u1	0.00
73) Anthracene-d10	17.463	188	1357307	37.545	ng/u1	0.00
81) Pyrene-d10	19.704	212	1812409	42.974	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.816	264	1550715	37.467	ng/u1	-0.02

Target Compounds Qvalue

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

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