

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
 Data File : BP017294.D
 Acq On : 22 Sep 2023 19:26
 Operator : MA/JU
 Sample : 04410-09
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHE1

Quant Time: Sep 22 22:32:36 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	7.934	152	139413	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	554620	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	330149	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	758966	20.000	ng/ul	0.00
79) Chrysene-d12	21.469	240	734573	20.000	ng/ul	0.00
88) Perylene-d12	23.992	264	826789	20.000	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	16028	4.722	ng/uL	0.00
4) Pyridine-d5	3.746	84	62498	6.587	ng/ul	0.00
7) Phenol-d5	7.093	99	89317	7.801	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	252240	36.504	ng/ul	0.00
11) 2-Chlorophenol-d4	7.457	132	271390	29.884	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	161038	18.090	ng/ul	-0.01
21) Nitrobenzene-d5	9.134	128	158796	39.604	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	169991	39.046	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	300154	36.491	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.922	131	250951	21.062	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	1064683	45.017	ng/ul	0.00
49) Acenaphthylene-d8	14.292	160	1125286	41.369	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	26988	6.550	ng/ul	0.00
60) Fluorene-d10	15.592	176	922022	45.283	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	142763	33.374	ng/ul	0.00
73) Anthracene-d10	17.463	188	1591442	47.279	ng/ul	0.00
81) Pyrene-d10	19.710	212	1970995	49.618	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264	1973991	49.589	ng/ul	0.00

Target Compounds Qvalue

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

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