

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
 Data File : BP017296.D
 Acq On : 22 Sep 2023 20:34
 Operator : MA/JU
 Sample : 04410-11
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJM0

Quant Time: Sep 22 22:34:59 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	139555	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	565336	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	338983	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	791470	20.000	ng/ul	0.00
79) Chrysene-d12	21.469	240	757733	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	846948	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	18051	5.313	ng/uL	0.00
4) Pyridine-d5	3.746	84	64420	6.782	ng/ul	0.00
7) Phenol-d5	7.093	99	84333	7.358	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	256773	37.122	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	261467	28.763	ng/ul	-0.01
15) 4-Methylphenol-d8	8.646	113	150888	16.933	ng/ul	-0.01
21) Nitrobenzene-d5	9.134	128	163892	40.100	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	169986	38.304	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	300701	35.865	ng/ul	0.00
31) 4-Chloroaniline-d4	10.922	131	187371	15.428	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	1102494	45.401	ng/ul	0.00
49) Acenaphthylene-d8	14.292	160	1180167	42.256	ng/ul	0.00
54) 4-Nitrophenol-d4	14.834	143	27413	6.480	ng/ul	0.00
60) Fluorene-d10	15.592	176	968095	46.307	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	143741	32.223	ng/ul	0.00
73) Anthracene-d10	17.463	188	1617771	46.087	ng/ul	0.00
81) Pyrene-d10	19.710	212	2046791	49.951	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1933439	47.414	ng/ul	-0.01

Target Compounds Qvalue

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

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