

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
 Data File : BP017295.D
 Acq On : 22 Sep 2023 20:00
 Operator : MA/JU
 Sample : 04410-10
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJL8

Quant Time: Sep 22 22:26:25 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	54464	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	221012	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.593	164	127492	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	291525	20.000	ng/ul	0.00
79) Chrysene-d12	21.469	240	281596	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	309047	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	15740	11.870	ng/uL	0.00
4) Pyridine-d5	3.746	84	68306	18.426	ng/ul	0.00
7) Phenol-d5	7.093	99	90889	20.319	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	95462	35.363	ng/ul	0.00
11) 2-Chlorophenol-d4	7.458	132	109706	30.923	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	83074	23.888	ng/ul	-0.01
21) Nitrobenzene-d5	9.134	128	54475	34.094	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	56200	32.394	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	102262	31.199	ng/ul	0.00
31) 4-Chloroaniline-d4	10.922	131	105516	22.223	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	326189	35.715	ng/ul	0.00
49) Acenaphthylene-d8	14.293	160	373374	35.545	ng/ul	0.00
54) 4-Nitrophenol-d4	14.840	143	21974	13.810	ng/ul	0.01
60) Fluorene-d10	15.592	176	318436	40.499	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	31529	19.189	ng/ul	0.00
73) Anthracene-d10	17.463	188	587054	45.405	ng/ul	0.00
81) Pyrene-d10	19.704	212	766715	50.349	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.816	264	751158	50.482	ng/ul	-0.02

Target Compounds Qvalue

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

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