

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017577.D
 Acq On : 05 Oct 2023 21:39
 Operator : MA/JU
 Sample : 04607-02
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJ35

Quant Time: Oct 06 02:03:46 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 04 16:33:47 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	88558	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.775	136	366933	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	235835	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.416	188	559934	20.000	ng/ul	-0.01
79) Chrysene-d12	21.551	240	594794	20.000	ng/ul	-0.01
88) Perylene-d12	24.133	264	676933	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	18596	8.395	ng/uL	0.00
4) Pyridine-d5	3.734	84	31601	5.450	ng/ul	0.00
7) Phenol-d5	7.093	99	60563	8.128	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	193601	42.582	ng/ul	0.00
11) 2-Chlorophenol-d4	7.458	132	190875	32.534	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	121053	20.294	ng/ul	-0.01
21) Nitrobenzene-d5	9.146	128	123328	43.688	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	131696	41.363	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	227702	38.576	ng/ul	0.00
31) 4-Chloroaniline-d4	10.946	131	301462	35.852	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	823812	44.783	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	904867	44.608	ng/ul	0.00
54) 4-Nitrophenol-d4	14.887	143	21177	6.574	ng/ul	0.01
60) Fluorene-d10	15.634	176	760359	48.730	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	122179	35.000	ng/ul	0.00
73) Anthracene-d10	17.516	188	1376820	53.121	ng/ul	-0.01
81) Pyrene-d10	19.775	212	1758052	53.192	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.963	264	1833623	53.674	ng/ul	-0.01

Target Compounds Qvalue

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

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