

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017605.D
 Acq On : 06 Oct 2023 17:04
 Operator : MA/JU
 Sample : 04624-09
 Misc :
 ALS Vial : 85 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Oct 06 23:14:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 06 05:55:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	94520	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	360413	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	226672	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.416	188	506727	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	520207	20.000	ng/ul	0.00
88) Perylene-d12	24.127	264	600683	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	11360	4.805	ng/uL	0.00
4) Pyridine-d5	3.728	84	79051	12.774	ng/ul	0.00
7) Phenol-d5	7.093	99	73696	9.267	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	211595	43.604	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	219064	34.984	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	135702	21.315	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	133448	48.128	ng/ul	0.00
24) 2-Nitrophenol-d4	9.858	143	140781	45.017	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	238026	41.055	ng/ul	0.00
31) 4-Chloroaniline-d4	10.940	131	294103	35.610	ng/ul	0.00
46) Dimethylphthalate-d6	14.040	166	798628	45.169	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	880977	45.186	ng/ul	0.00
54) 4-Nitrophenol-d4	14.887	143	18348	5.926	ng/ul	0.02
60) Fluorene-d10	15.628	176	696457	46.439	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200	195	0.062	ng/ul	0.09
73) Anthracene-d10	17.516	188	1154325	49.213	ng/ul	0.00
81) Pyrene-d10	19.769	212	1420035	49.126	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264	1533244	50.578	ng/ul	0.00

Target Compounds	Qvalue

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

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