

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017607.D
 Acq On : 06 Oct 2023 18:12
 Operator : MA/JU
 Sample : 04624-11
 Misc :
 ALS Vial : 87 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Oct 06 23:15:04 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	98880	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	417180	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	262628	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.416	188	620727	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	674694	20.000	ng/ul	0.00
88) Perylene-d12	24.127	264	782735	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	11104	4.490	ng/uL	0.00
4) Pyridine-d5	3.728	84	65293	10.086	ng/ul	0.00
7) Phenol-d5	7.093	99	65945	7.927	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	190120	37.451	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	197141	30.095	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	117604	17.658	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	122535	38.179	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143	131220	36.250	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	234349	34.921	ng/ul	0.00
31) 4-Chloroaniline-d4	10.940	131	252989	26.463	ng/ul	0.00
46) Dimethylphthalate-d6	14.040	166	841764	41.090	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	892844	39.525	ng/ul	0.00
54) 4-Nitrophenol-d4	14.881	143	21943	6.117	ng/ul	0.01
60) Fluorene-d10	15.628	176	754940	43.447	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	120698	31.190	ng/ul	0.00
73) Anthracene-d10	17.516	188	1273868	44.335	ng/ul	0.00
81) Pyrene-d10	19.769	212	1588933	42.382	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264	1713641	43.381	ng/ul	0.00

Target Compounds Qvalue

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

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