

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032824\
 Data File : BP020017.D
 Acq On : 28 Mar 2024 19:30
 Operator : MA/JU
 Sample : P1840-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Mar 28 23:09:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 28 14:08:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	127739	20.000	ng/u1	0.00
20) Naphthalene-d8	10.357	136	504038	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.245	164	375019	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.063	188	812458	20.000	ng/u1	-0.01
79) Chrysene-d12	21.539	240	717980	20.000	ng/u1	0.00
88) Perylene-d12	24.833	264	773169	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	71	0.024	ng/uL	0.01
4) Pyridine-d5	3.634	84	478	0.054	ng/u1	0.00
7) Phenol-d5	6.793	99	18	0.002	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	60	0.009	ng/u1	0.00
11) 2-Chlorophenol-d4	7.157	132	9	0.001	ng/u1	0.00
15) 4-Methylphenol-d8	8.293	113	17	0.002	ng/u1	-0.01
21) Nitrobenzene-d5	8.740	128	13	0.003	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.446	143	18	0.004	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	9.987	165	51	0.006	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.516	131	71	0.007	ng/u1	0.00
46) Dimethylphthalate-d6	13.651	166	4	0.000	ng/u1	0.00
49) Acenaphthylene-d8	13.939	160	43	0.001	ng/u1	0.00
54) 4-Nitrophenol-d4	14.481	143	20	0.005	ng/u1	0.00
60) Fluorene-d10	15.251	176	13	0.001	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	5	0.001	ng/u1	-0.01
73) Anthracene-d10	17.063	188	812458	22.591	ng/u1	-0.11
81) Pyrene-d10	19.569	212	10	0.000	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.609	264	317	0.009	ng/u1	0.00

Target Compounds Qvalue

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

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