

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
 Data File : BP017487.D
 Acq On : 02 Oct 2023 20:50
 Operator : MA/JU
 Sample : 04571-17
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBK95

Quant Time: Oct 02 23:16:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 30 04:02:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.945	152	124885	20.000	ng/u1	0.00
20) Naphthalene-d8	10.781	136	509903	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.639	164	304568	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.427	188	691220	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	661029	20.000	ng/u1	0.00
88) Perylene-d12	24.127	264	734761	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	17395	5.721	ng/uL	0.00
4) Pyridine-d5	3.734	84	111969	13.173	ng/u1	0.00
7) Phenol-d5	7.098	99	98403	9.594	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	260689	42.115	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	274569	33.752	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	187231	23.480	ng/u1	0.00
21) Nitrobenzene-d5	9.157	128	169911	46.093	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	180311	45.048	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	326350	43.156	ng/u1	0.00
31) 4-Chloroaniline-d4	10.951	131	310963	28.387	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	1110719	50.908	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	1228733	48.966	ng/u1	0.00
54) 4-Nitrophenol-d4	14.880	143	23358	6.145	ng/u1	0.00
60) Fluorene-d10	15.645	176	957880	50.996	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.780	200	150666	38.673	ng/u1	0.00
73) Anthracene-d10	17.527	188	1625099	53.011	ng/u1	0.00
81) Pyrene-d10	19.780	212	2005706	56.109	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.956	264	1923877	54.383	ng/u1	0.00

Target Compounds Qvalue

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

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