

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021225\
 Data File : BP024053.D
 Acq On : 12 Feb 2025 17:03
 Operator : RC/JU
 Sample : P5323-18
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YE641

Quant Time: Feb 14 11:38:52 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 13 12:56:03 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.758	152	527112	20.000	ng/u1	0.00
20) Naphthalene-d8	10.534	136	2031702	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.375	164	1146834	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.163	188	2006821	20.000	ng/u1	0.00
79) Chrysene-d12	21.604	240	2005562	20.000	ng/u1	0.00
88) Perylene-d12	24.945	264	2186953	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	105	0.008	ng/uL	0.01
4) Pyridine-d5	3.693	84	168	0.005	ng/u1	0.00
7) Phenol-d5	6.952	99	180	0.004	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.099	67	3669	0.149	ng/u1	0.00
11) 2-Chlorophenol-d4	7.305	132	26	0.001	ng/u1	0.00
15) 4-Methylphenol-d8	8.475	113	245	0.006	ng/u1	0.01
21) Nitrobenzene-d5	8.893	128	23	0.001	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.634	143	5	0.000	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.157	165	138	0.004	ng/u1	0.00
31) 4-Chloroaniline-d4	10.646	131	36614	0.761	ng/u1	-0.02
46) Dimethylphthalate-d6	13.787	166	301	0.004	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	377	0.004	ng/u1	0.00
54) 4-Nitrophenol-d4	14.598	143	2536	0.149	ng/u1	0.00
60) Fluorene-d10	15.381	176	277	0.004	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	974	0.079	ng/u1	-0.02
73) Anthracene-d10	0.000	188	0d	0.000	ng/u1	
81) Pyrene-d10	19.639	212	2092	0.019	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.727	264	1192	0.010	ng/u1	0.00

Target Compounds	Qvalue

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

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