

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
 Data File : BP020905.D
 Acq On : 11 Jul 2024 15:56
 Operator : MA/JU
 Sample : P3087-19
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4BR4

Quant Time: Jul 11 23:23:38 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 10 18:38:19 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	220208	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.481	136	956916	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.357	164	637837	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.157	188	1179312	20.000	ng/u1	-0.02
79) Chrysene-d12	21.627	240	860722	20.000	ng/u1	0.00
88) Perylene-d12	24.962	264	1054907	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	19617	4.054	ng/uL	0.00
4) Pyridine-d5	3.664	84	263548	18.812	ng/u1	0.00
7) Phenol-d5	6.911	99	417646	23.263	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.058	67	272382	25.035	ng/u1	0.00
11) 2-Chlorophenol-d4	7.263	132	359954	24.335	ng/u1	0.00
15) 4-Methylphenol-d8	8.422	113	359525	24.112	ng/u1	-0.01
21) Nitrobenzene-d5	8.869	128	182332	24.532	ng/u1	0.00
24) 2-Nitrophenol-d4	9.575	143	208903	24.557	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.122	165	353712	22.994	ng/u1	0.00
31) 4-Chloroaniline-d4	10.628	131	475752	23.346	ng/u1	-0.01
46) Dimethylphthalate-d6	13.757	166	1154444	24.898	ng/u1	-0.01
49) Acenaphthylene-d8	14.045	160	1280243	24.590	ng/u1	0.00
54) 4-Nitrophenol-d4	14.598	143	115144	17.454	ng/u1	-0.01
60) Fluorene-d10	15.357	176	944386	24.827	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.504	200	127219	17.829	ng/u1	0.00
73) Anthracene-d10	17.257	188	1286719	24.565	ng/u1	-0.02
81) Pyrene-d10	19.651	212	1262936	23.733	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.727	264	1304748	25.078	ng/u1	-0.03

Target Compounds	Qvalue

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

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