

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
 Data File : BP020927.D
 Acq On : 12 Jul 2024 09:18
 Operator : MA/JU
 Sample : P3087-15
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4BR0

Quant Time: Jul 12 23:07:20 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	7.722	152	205953	20.000	ng/u1	0.00
20) Naphthalene-d8	10.487	136	780197	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.345	164	475828	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.163	188	1039969	20.000	ng/u1	-0.02
79) Chrysene-d12	21.616	240	1096989	20.000	ng/u1	0.00
88) Perylene-d12	24.974	264	1329335	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	23285	5.145	ng/uL	0.00
4) Pyridine-d5	3.670	84	337701	25.773	ng/u1	0.00
7) Phenol-d5	6.916	99	542461	32.307	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.069	67	335544	32.975	ng/u1	0.00
11) 2-Chlorophenol-d4	7.263	132	462000	33.396	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	442236	31.712	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	218670	36.085	ng/u1	0.00
24) 2-Nitrophenol-d4	9.587	143	252916	36.465	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	427758	34.107	ng/u1	0.00
31) 4-Chloroaniline-d4	10.634	131	586000	35.270	ng/u1	0.00
46) Dimethylphthalate-d6	13.757	166	1399247	40.452	ng/u1	-0.01
49) Acenaphthylene-d8	14.040	160	1540613	39.666	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.610	143	212911	43.262	ng/u1	0.00
60) Fluorene-d10	15.363	176	1205972	42.498	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.510	200	205667	32.685	ng/u1	0.00
73) Anthracene-d10	17.269	188	1928730	41.756	ng/u1	0.00
81) Pyrene-d10	19.651	212	2416572	35.632	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.745	264	2869055	43.761	ng/u1	-0.01

Target Compounds	Qvalue

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

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