

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
 Data File : BP020922.D
 Acq On : 12 Jul 2024 05:01
 Operator : MA/JU
 Sample : P3087-13
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4BP1

Quant Time: Jul 12 05:29:56 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	245233	20.000	ng/u1	0.00
20) Naphthalene-d8	10.493	136	1012069	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.351	164	685287	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.163	188	1507611	20.000	ng/u1	-0.02
79) Chrysene-d12	21.622	240	1471056	20.000	ng/u1	0.00
88) Perylene-d12	24.968	264	1681910	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	14633	2.715	ng/uL	0.00
4) Pyridine-d5	3.675	84	214097	13.723	ng/u1	0.00
7) Phenol-d5	6.922	99	340599	17.036	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.063	67	212220	17.515	ng/u1	0.00
11) 2-Chlorophenol-d4	7.269	132	290040	17.607	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	295011	17.766	ng/u1	0.00
21) Nitrobenzene-d5	8.869	128	143951	18.312	ng/u1	0.00
24) 2-Nitrophenol-d4	9.581	143	164277	18.259	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	297214	18.269	ng/u1	0.01
31) 4-Chloroaniline-d4	10.634	131	398862	18.506	ng/u1	0.00
46) Dimethylphthalate-d6	13.763	166	1029724	20.670	ng/u1	0.00
49) Acenaphthylene-d8	14.045	160	1140005	20.380	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	156423	22.069	ng/u1	0.00
60) Fluorene-d10	15.363	176	899125	22.000	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.510	200	135841	14.892	ng/u1	0.00
73) Anthracene-d10	17.263	188	1467216	21.912	ng/u1	-0.01
81) Pyrene-d10	19.651	212	1801703	19.811	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.745	264	1839606	22.177	ng/u1	-0.01

Target Compounds	Qvalue

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

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