

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013025\
 Data File : BP023808.D
 Acq On : 30 Jan 2025 21:49
 Operator : RC/JU
 Sample : Q1192-05
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A44W0

Quant Time: Jan 31 08:15:02 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 31 08:09:38 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	419803	20.000	ng/u1	0.02
20) Naphthalene-d8	10.557	136	1794149	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.398	164	1202974	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.187	188	2454512	20.000	ng/u1	-0.02
79) Chrysene-d12	21.627	240	2422737	20.000	ng/u1	-0.02
88) Perylene-d12	24.992	264	2640768	20.000	ng/u1	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	49830	4.932	ng/uL	0.01
4) Pyridine-d5	3.711	84	220214	7.553	ng/u1	0.02
7) Phenol-d5	6.964	99	303234	8.162	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.116	67	694807	35.333	ng/u1	0.02
11) 2-Chlorophenol-d4	7.328	132	888882	30.468	ng/u1	0.02
15) 4-Methylphenol-d8	8.487	113	601274	19.812	ng/u1	0.02
21) Nitrobenzene-d5	8.928	128	501183	34.915	ng/u1	0.02
24) 2-Nitrophenol-d4	9.646	143	523526	33.801	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.187	165	921189	32.420	ng/u1	0.01
31) 4-Chloroaniline-d4	10.687	131	1217044	28.641	ng/u1	0.00
46) Dimethylphthalate-d6	13.804	166	3523941	39.273	ng/u1	0.00
49) Acenaphthylene-d8	14.087	160	3695267	36.514	ng/u1	0.00
54) 4-Nitrophenol-d4	14.604	143	143190	8.035	ng/u1	-0.02
60) Fluorene-d10	15.404	176	3017638	39.935	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.516	200	513569	33.916	ng/u1	-0.02
73) Anthracene-d10	17.292	188	5460575	45.310	ng/u1	-0.02
81) Pyrene-d10	19.657	212	6320827	48.691	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.768	264	6418928	46.587	ng/u1	-0.02

Target Compounds Qvalue

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

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