

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013025\
 Data File : BP023810.D
 Acq On : 30 Jan 2025 23:10
 Operator : RC/JU
 Sample : Q1192-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A44W2

Quant Time: Jan 31 08:15:31 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	442151	20.000	ng/u1	0.02
20) Naphthalene-d8	10.557	136	1842468	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.404	164	1207184	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.192	188	2433204	20.000	ng/u1	-0.02
79) Chrysene-d12	21.628	240	2519667	20.000	ng/u1	-0.02
88) Perylene-d12	25.016	264	2748907	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	51333	4.824	ng/uL	0.01
4) Pyridine-d5	3.717	84	174864	5.695	ng/u1	0.02
7) Phenol-d5	6.964	99	294446	7.525	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.117	67	619235	29.898	ng/u1	0.02
11) 2-Chlorophenol-d4	7.328	132	799660	26.025	ng/u1	0.02
15) 4-Methylphenol-d8	8.481	113	544009	17.019	ng/u1	0.01
21) Nitrobenzene-d5	8.922	128	434301	29.462	ng/u1	0.01
24) 2-Nitrophenol-d4	9.640	143	450520	28.324	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.187	165	807838	27.685	ng/u1	0.01
31) 4-Chloroaniline-d4	10.687	131	1009845	23.142	ng/u1	0.00
46) Dimethylphthalate-d6	13.810	166	2992690	33.236	ng/u1	0.00
49) Acenaphthylene-d8	14.093	160	3180079	31.313	ng/u1	0.00
54) 4-Nitrophenol-d4	14.616	143	124410	6.957	ng/u1	0.00
60) Fluorene-d10	15.410	176	2581664	34.046	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.516	200	400083	26.653	ng/u1	-0.02
73) Anthracene-d10	17.292	188	4499115	37.659	ng/u1	-0.02
81) Pyrene-d10	19.663	212	5347669	39.610	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.786	264	5604809	39.078	ng/u1	0.00

Target Compounds	Qvalue

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

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