

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013025\
 Data File : BP023807.D
 Acq On : 30 Jan 2025 21:08
 Operator : RC/JU
 Sample : Q1192-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A44T9

Quant Time: Jan 31 08:14:48 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	468611	20.000	ng/u1	0.02
20) Naphthalene-d8	10.557	136	1976543	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.398	164	1326150	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.186	188	2708667	20.000	ng/u1	-0.02
79) Chrysene-d12	21.627	240	2727703	20.000	ng/u1	-0.02
88) Perylene-d12	24.998	264	2925363	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	62896	5.577	ng/uL	0.00
4) Pyridine-d5	3.711	84	353342	10.857	ng/u1	0.02
7) Phenol-d5	6.969	99	315448	7.607	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.116	67	703573	32.052	ng/u1	0.02
11) 2-Chlorophenol-d4	7.328	132	901081	27.669	ng/u1	0.02
15) 4-Methylphenol-d8	8.487	113	611024	18.037	ng/u1	0.02
21) Nitrobenzene-d5	8.928	128	510311	32.270	ng/u1	0.02
24) 2-Nitrophenol-d4	9.646	143	531127	31.127	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.187	165	920619	29.410	ng/u1	0.01
31) 4-Chloroaniline-d4	10.693	131	1376449	29.403	ng/u1	0.01
46) Dimethylphthalate-d6	13.804	166	3510916	35.493	ng/u1	0.00
49) Acenaphthylene-d8	14.092	160	3752577	33.636	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	139177	7.085	ng/u1	-0.01
60) Fluorene-d10	15.398	176	3067254	36.821	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.516	200	478261	28.621	ng/u1	-0.02
73) Anthracene-d10	17.292	188	5242574	39.419	ng/u1	-0.02
81) Pyrene-d10	19.657	212	6019398	41.185	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.763	264	6304473	41.305	ng/u1	-0.02

Target Compounds	Qvalue

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

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