

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
 Data File : BP023849.D
 Acq On : 01 Feb 2025 07:41
 Operator : RC/JU
 Sample : Q1224-16
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A6323

Quant Time: Feb 03 08:15:47 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 31 11:40:55 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	492374	20.000	ng/u1	0.02
20) Naphthalene-d8	10.558	136	2080522	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.404	164	1327636	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.192	188	2696092	20.000	ng/u1	-0.02
79) Chrysene-d12	21.633	240	2834922	20.000	ng/u1	-0.01
88) Perylene-d12	25.010	264	3082261	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	51193	4.320	ng/uL	0.00
4) Pyridine-d5	3.711	84	574150	16.790	ng/u1	0.02
7) Phenol-d5	6.969	99	929524	21.332	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.117	67	527546	22.873	ng/u1	0.02
11) 2-Chlorophenol-d4	7.322	132	784693	22.933	ng/u1	0.01
15) 4-Methylphenol-d8	8.487	113	757523	21.282	ng/u1	0.02
21) Nitrobenzene-d5	8.928	128	371429	22.314	ng/u1	0.02
24) 2-Nitrophenol-d4	9.646	143	400507	22.299	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.187	165	698823	21.209	ng/u1	0.01
31) 4-Chloroaniline-d4	10.687	131	962332	19.529	ng/u1	0.00
46) Dimethylphthalate-d6	13.810	166	2117180	21.379	ng/u1	0.00
49) Acenaphthylene-d8	14.099	160	2139008	19.151	ng/u1	0.00
54) 4-Nitrophenol-d4	14.622	143	321644	16.355	ng/u1	0.00
60) Fluorene-d10	15.410	176	1493471	17.909	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.522	200	186401	11.207	ng/u1	-0.01
73) Anthracene-d10	17.298	188	2192787	16.565	ng/u1	-0.02
81) Pyrene-d10	19.663	212	2387181	15.715	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.769	264	2056487	12.788	ng/u1	-0.02

Target Compounds Qvalue

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

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