

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
 Data File : BP023859.D
 Acq On : 01 Feb 2025 15:29
 Operator : RC/JU
 Sample : Q1248-09
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A6342

Quant Time: Feb 03 08:33:51 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	390682	20.000	ng/u1	0.02
20) Naphthalene-d8	10.552	136	1672304	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.398	164	1076503	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.192	188	2118020	20.000	ng/u1	-0.02
79) Chrysene-d12	21.633	240	2071934	20.000	ng/u1	-0.01
88) Perylene-d12	24.998	264	2168613	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	56914	6.053	ng/uL	0.01
4) Pyridine-d5	3.705	84	791559	29.174	ng/u1	0.01
7) Phenol-d5	6.964	99	1102016	31.874	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.111	67	610564	33.363	ng/u1	0.01
11) 2-Chlorophenol-d4	7.322	132	902211	33.230	ng/u1	0.01
15) 4-Methylphenol-d8	8.481	113	914713	32.387	ng/u1	0.01
21) Nitrobenzene-d5	8.922	128	431024	32.215	ng/u1	0.01
24) 2-Nitrophenol-d4	9.634	143	456735	31.637	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.181	165	834689	31.516	ng/u1	0.00
31) 4-Chloroaniline-d4	10.681	131	1160182	29.292	ng/u1	0.00
46) Dimethylphthalate-d6	13.804	166	2659348	33.119	ng/u1	0.00
49) Acenaphthylene-d8	14.093	160	2724192	30.081	ng/u1	0.00
54) 4-Nitrophenol-d4	14.616	143	422338	26.485	ng/u1	0.00
60) Fluorene-d10	15.410	176	1949778	28.835	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.516	200	219169	16.773	ng/u1	-0.02
73) Anthracene-d10	17.292	188	2809134	27.012	ng/u1	-0.02
81) Pyrene-d10	19.669	212	2878234	25.926	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.763	264	2341761	20.696	ng/u1	-0.02

Target Compounds Qvalue

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

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