

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011725\
 Data File : BP023631.D
 Acq On : 17 Jan 2025 16:40
 Operator : RC/JU
 Sample : Q1086-13
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AF1

Quant Time: Jan 17 17:15:44 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 14 16:27:27 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.811	152	538303	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	2284064	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.428	164	1466047	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.222	188	2869271	20.000	ng/u1	-0.02
79) Chrysene-d12	21.692	240	2747371	20.000	ng/u1	-0.02
88) Perylene-d12	25.127	264	2909679	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	72217	4.947	ng/uL	0.00
4) Pyridine-d5	3.723	84	282755	6.947	ng/u1	0.00
7) Phenol-d5	6.969	99	399491	8.620	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	865717	31.888	ng/u1	0.00
11) 2-Chlorophenol-d4	7.340	132	1125146	30.742	ng/u1	0.00
15) 4-Methylphenol-d8	8.493	113	722437	20.097	ng/u1	0.00
21) Nitrobenzene-d5	8.946	128	556809	31.939	ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	495023	43.259	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	1284790	36.913	ng/u1	0.00
31) 4-Chloroaniline-d4	10.710	131	1118057	20.121	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	3959153	34.700	ng/u1	0.00
49) Acenaphthylene-d8	14.116	160	4424999	32.602	ng/u1	0.00
54) 4-Nitrophenol-d4	14.604	143	153350	8.086	ng/u1	0.00
60) Fluorene-d10	15.434	176	3405666	34.490	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.540	200	296061	31.712	ng/u1	-0.01
73) Anthracene-d10	17.322	188	5465990	35.290	ng/u1	-0.02
81) Pyrene-d10	19.704	212	6313404	41.481	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.880	264	5903325	37.333	ng/u1	-0.03

Target Compounds	Qvalue

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

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