

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012425\
 Data File : BP023721.D
 Acq On : 24 Jan 2025 19:21
 Operator : RC/JU
 Sample : Q1088-10
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B09

Quant Time: Jan 24 22:26:21 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012225.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 23 00:03:00 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	503030	20.000	ng/u1	0.00
20) Naphthalene-d8	10.563	136	2056820	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.416	164	1277003	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.210	188	2570214	20.000	ng/u1	0.00
79) Chrysene-d12	21.651	240	2649621	20.000	ng/u1	0.00
88) Perylene-d12	25.045	264	2819290	20.000	ng/u1	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	67839	5.432	ng/uL	0.00
4) Pyridine-d5	3.711	84	201209	5.583	ng/u1	0.00
7) Phenol-d5	6.952	99	333420	7.833	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	865463	35.998	ng/u1	0.00
11) 2-Chlorophenol-d4	7.328	132	1017047	31.430	ng/u1	0.00
15) 4-Methylphenol-d8	8.481	113	658926	19.687	ng/u1	0.00
21) Nitrobenzene-d5	8.934	128	555615	35.854	ng/u1	0.00
24) 2-Nitrophenol-d4	9.652	143	516580	36.360	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.181	165	1020980	34.189	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.693	131	1257669	26.334	ng/u1	0.00
46) Dimethylphthalate-d6	13.816	166	3435462	37.408	ng/u1	0.00
49) Acenaphthylene-d8	14.104	160	3966143	37.314	ng/u1	0.00
54) 4-Nitrophenol-d4	14.592	143	115494	6.259	ng/u1	-0.01
60) Fluorene-d10	15.422	176	3105904	39.508	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.528	200	342616	27.069	ng/u1	0.00
73) Anthracene-d10	17.310	188	5256808	42.602	ng/u1	0.00
81) Pyrene-d10	19.680	212	6284078	46.609	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.810	264	6268363	44.420	ng/u1	0.00

Target Compounds	Qvalue

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

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