

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
 Data File : BP023864.D
 Acq On : 01 Feb 2025 19:32
 Operator : RC/JU
 Sample : PB166436BL
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK436

Quant Time: Feb 03 08:24:10 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	386141	20.000	ng/u1	0.02
20) Naphthalene-d8	10.557	136	1578343	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.398	164	1028600	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.192	188	2119117	20.000	ng/u1	-0.02
79) Chrysene-d12	21.633	240	2143537	20.000	ng/u1	-0.01
88) Perylene-d12	25.004	264	2315186	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	69231	7.449	ng/uL	0.00
4) Pyridine-d5	3.705	84	1005695	37.502	ng/u1	0.01
7) Phenol-d5	6.963	99	1278169	37.404	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.116	67	641221	35.450	ng/u1	0.02
11) 2-Chlorophenol-d4	7.322	132	1028484	38.327	ng/u1	0.01
15) 4-Methylphenol-d8	8.487	113	1014641	36.347	ng/u1	0.02
21) Nitrobenzene-d5	8.922	128	455704	36.088	ng/u1	0.01
24) 2-Nitrophenol-d4	9.640	143	474221	34.804	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.181	165	926605	37.070	ng/u1	0.00
31) 4-Chloroaniline-d4	10.687	131	1333475	35.671	ng/u1	0.00
46) Dimethylphthalate-d6	13.804	166	2697098	35.153	ng/u1	0.00
49) Acenaphthylene-d8	14.087	160	3156285	36.475	ng/u1	0.00
54) 4-Nitrophenol-d4	14.616	143	471018	30.913	ng/u1	0.00
60) Fluorene-d10	15.404	176	2438224	37.737	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.516	200	274684	21.011	ng/u1	-0.02
73) Anthracene-d10	17.286	188	3932621	37.796	ng/u1	-0.03
81) Pyrene-d10	19.657	212	4528534	39.428	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.763	264	4610943	38.171	ng/u1	-0.02

Target Compounds Qvalue

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

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