

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
 Data File : BP023836.D
 Acq On : 31 Jan 2025 22:48
 Operator : RC/JU
 Sample : PB166393BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK393

Quant Time: Feb 03 07:37:16 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	442041	20.000	ng/u1	0.02
20) Naphthalene-d8	10.558	136	1904768	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.404	164	1316077	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.204	188	2767373	20.000	ng/u1	0.00
79) Chrysene-d12	21.639	240	2811171	20.000	ng/u1	0.00
88) Perylene-d12	25.027	264	3042054	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	63027	5.924	ng/uL	0.00
4) Pyridine-d5	3.705	84	933424	30.405	ng/u1	0.01
7) Phenol-d5	6.969	99	1214788	31.053	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.111	67	613327	29.620	ng/u1	0.01
11) 2-Chlorophenol-d4	7.322	132	965135	31.418	ng/u1	0.01
15) 4-Methylphenol-d8	8.487	113	980282	30.676	ng/u1	0.02
21) Nitrobenzene-d5	8.922	128	444929	29.196	ng/u1	0.01
24) 2-Nitrophenol-d4	9.640	143	462166	28.106	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.187	165	910077	30.169	ng/u1	0.01
31) 4-Chloroaniline-d4	10.693	131	1301903	28.859	ng/u1	0.01
46) Dimethylphthalate-d6	13.810	166	2919031	29.735	ng/u1	0.00
49) Acenaphthylene-d8	14.099	160	3228481	29.160	ng/u1	0.00
54) 4-Nitrophenol-d4	14.628	143	497639	25.526	ng/u1	0.00
60) Fluorene-d10	15.410	176	2491047	30.133	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.522	200	321749	18.846	ng/u1	-0.01
73) Anthracene-d10	17.304	188	4148863	30.534	ng/u1	-0.01
81) Pyrene-d10	19.669	212	4773716	31.692	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.798	264	4896768	30.852	ng/u1	0.01

Target Compounds Qvalue

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

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