

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
 Data File : BP023852.D
 Acq On : 01 Feb 2025 10:38
 Operator : RC/JU
 Sample : PB166436BL
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK436

Quant Time: Feb 03 08:22:19 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	7.757	152	476755	20.000	ng/ul	0.00
20) Naphthalene-d8	10.533	136	1973526	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.398	164	1311821	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	2770162	20.000	ng/ul	0.00
79) Chrysene-d12	21.633	240	2832875	20.000	ng/ul	-0.01
88) Perylene-d12	25.021	264	3031336	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	66667	5.810	ng/uL	-0.01
4) Pyridine-d5	3.687	84	993140	29.995	ng/ul	0.00
7) Phenol-d5	6.940	99	1261855	29.908	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.092	67	626667	28.061	ng/ul	0.00
11) 2-Chlorophenol-d4	7.298	132	1009512	30.469	ng/ul	-0.01
15) 4-Methylphenol-d8	8.463	113	989185	28.701	ng/ul	0.00
21) Nitrobenzene-d5	8.898	128	445873	28.239	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.622	143	470183	27.597	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.163	165	918868	29.399	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.669	131	1307527	27.973	ng/ul	-0.01
46) Dimethylphthalate-d6	13.798	166	2727289	27.872	ng/ul	0.00
49) Acenaphthylene-d8	14.086	160	3129394	28.356	ng/ul	0.00
54) 4-Nitrophenol-d4	14.621	143	496835	25.568	ng/ul	0.00
60) Fluorene-d10	15.398	176	2410587	29.254	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.521	200	315248	18.447	ng/ul	-0.01
73) Anthracene-d10	17.304	188	3968825	29.179	ng/ul	-0.01
81) Pyrene-d10	19.668	212	4565658	30.079	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.780	264	4608413	29.137	ng/ul	0.00

Target Compounds Qvalue

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

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