

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
 Data File : BP023938.D
 Acq On : 05 Feb 2025 16:24
 Operator : RC/JU
 Sample : PB166504BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK504

Quant Time: Feb 05 17:52:26 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 03 11:19:36 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	362736	20.000	ng/u1	0.00
20) Naphthalene-d8	10.546	136	1557033	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.387	164	1010979	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.175	188	2052543	20.000	ng/u1	-0.01
79) Chrysene-d12	21.604	240	1980266	20.000	ng/u1	-0.04
88) Perylene-d12	24.962	264	2084430	20.000	ng/u1	-0.05

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	53876	6.171	ng/uL	0.00
4) Pyridine-d5	3.699	84	808025	32.075	ng/u1	0.00
7) Phenol-d5	6.957	99	1021906	31.834	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	534641	31.465	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	811355	32.186	ng/u1	0.00
15) 4-Methylphenol-d8	8.475	113	809187	30.858	ng/u1	0.00
21) Nitrobenzene-d5	8.916	128	370965	29.779	ng/u1	0.00
24) 2-Nitrophenol-d4	9.634	143	382009	28.420	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.175	165	746125	30.258	ng/u1	0.00
31) 4-Chloroaniline-d4	10.675	131	1097891	29.771	ng/u1	0.00
46) Dimethylphthalate-d6	13.792	166	2221032	29.453	ng/u1	0.00
49) Acenaphthylene-d8	14.081	160	2615404	30.751	ng/u1	0.00
54) 4-Nitrophenol-d4	14.598	143	358958	23.969	ng/u1	-0.01
60) Fluorene-d10	15.386	176	2000629	31.504	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.504	200	220901	17.445	ng/u1	0.00
73) Anthracene-d10	17.275	188	3185318	31.607	ng/u1	-0.01
81) Pyrene-d10	19.639	212	3651347	34.412	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.739	264	3517809	32.346	ng/u1	-0.04

Target Compounds	Qvalue

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

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