

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021225\
 Data File : BP024048.D
 Acq On : 12 Feb 2025 13:41
 Operator : RC/JU
 Sample : PB166649BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK649

Quant Time: Feb 13 17:12:20 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 13 12:56:03 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.740	152	336121	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.516	136	1466205	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.375	164	923343	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.180	188	1840054	20.000	ng/u1	0.01
79) Chrysene-d12	21.621	240	1687042	20.000	ng/u1	0.02
88) Perylene-d12	24.986	264	1659048	20.000	ng/u1	0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	56090	6.933	ng/uL	-0.02
4) Pyridine-d5	3.675	84	819213	35.094	ng/u1	-0.02
7) Phenol-d5	6.916	99	1060708	35.659	ng/u1	-0.03
9) Bis-(2-Chloroethyl)eth...	7.075	67	546391	34.703	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.281	132	863568	36.970	ng/u1	-0.02
15) 4-Methylphenol-d8	8.440	113	850178	34.988	ng/u1	-0.02
21) Nitrobenzene-d5	8.881	128	395651	33.728	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.599	143	405162	32.009	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.140	165	777886	33.500	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.646	131	1122365	32.320	ng/u1	-0.02
46) Dimethylphthalate-d6	13.781	166	2234630	32.446	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	2692975	34.668	ng/u1	0.00
54) 4-Nitrophenol-d4	14.592	143	296316	21.664	ng/u1	0.00
60) Fluorene-d10	15.386	176	2027859	34.964	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.504	200	156884	13.820	ng/u1	0.00
73) Anthracene-d10	17.286	188	3185009	35.253	ng/u1	0.02
81) Pyrene-d10	19.657	212	3518925	38.928	ng/u1	0.02
92) Benzo(a)pyrene-d12	24.762	264	3129421	36.153	ng/u1	0.04

Target Compounds Qvalue

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

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