

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
 Data File : BP021105.D
 Acq On : 23 Jul 2024 19:22
 Operator : MA/JU
 Sample : PB162098BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK098

Quant Time: Jul 24 04:12:45 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 24 00:18:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	150446	20.000	ng/u1	0.00
20) Naphthalene-d8	10.440	136	599755	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.304	164	391044	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.128	188	864840	20.000	ng/u1	0.00
79) Chrysene-d12	21.563	240	895672	20.000	ng/u1	-0.04
88) Perylene-d12	24.880	264	1090785	20.000	ng/u1	-0.06
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	20448	6.185	ng/uL	0.00
4) Pyridine-d5	3.611	84	259817	27.145	ng/u1	0.00
7) Phenol-d5	6.875	99	355893	29.016	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.010	67	209719	28.214	ng/u1	0.00
11) 2-Chlorophenol-d4	7.210	132	294157	29.108	ng/u1	-0.01
15) 4-Methylphenol-d8	8.393	113	283289	27.809	ng/u1	0.00
21) Nitrobenzene-d5	8.822	128	141157	30.302	ng/u1	0.00
24) 2-Nitrophenol-d4	9.534	143	166479	31.224	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	280573	29.102	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.587	131	336779	26.368	ng/u1	-0.01
46) Dimethylphthalate-d6	13.710	166	867463	30.516	ng/u1	-0.01
49) Acenaphthylene-d8	13.998	160	970016	30.390	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.598	143	131273	32.457	ng/u1	0.00
60) Fluorene-d10	15.322	176	744348	31.917	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.486	200	121018	23.127	ng/u1	-0.01
73) Anthracene-d10	17.222	188	1227663	31.960	ng/u1	-0.01
81) Pyrene-d10	19.604	212	1594237	28.790	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.651	264	1690209	31.418	ng/u1	-0.05

Target Compounds Qvalue

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

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