

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021641.D
 Acq On : 26 Aug 2024 16:38
 Operator : RC/JU
 Sample : PB162962BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK962

Quant Time: Aug 27 00:28:42 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 23 17:51:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	298422	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	1212373	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.439	164	811372	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	1793383	20.000	ng/u1	0.02
79) Chrysene-d12	21.710	240	1847559	20.000	ng/u1	0.00
88) Perylene-d12	25.139	264	2113199	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	44696	4.972	ng/uL	0.00
4) Pyridine-d5	3.687	84	604929	23.267	ng/u1	0.00
7) Phenol-d5	6.958	99	737170	23.147	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.122	67	423109	24.412	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.328	132	479447	23.318	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	556887	22.629	ng/u1	-0.01
21) Nitrobenzene-d5	8.940	128	239941	24.599	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	233832	23.617	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	442053	22.657	ng/u1	0.00
31) 4-Chloroaniline-d4	10.704	131	633516	22.318	ng/u1	0.00
46) Dimethylphthalate-d6	13.845	166	1600771	25.784	ng/u1	0.00
49) Acenaphthylene-d8	14.134	160	1735641	24.900	ng/u1	0.00
54) 4-Nitrophenol-d4	14.634	143	258278	21.866	ng/u1	0.00
60) Fluorene-d10	15.451	176	1349833	25.861	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	184731	18.728	ng/u1	0.00
73) Anthracene-d10	17.357	188	2182445	25.642	ng/u1	0.01
81) Pyrene-d10	19.727	212	2687013	25.442	ng/u1	0.01
92) Benzo(a)pyrene-d12	24.892	264	2771579	24.596	ng/u1	-0.01

Target Compounds Qvalue

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

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