

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080824\
 Data File : BG062352.D
 Acq On : 9 Aug 2024 9:08
 Operator : MA/JU
 Sample : PB162597BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK597

Quant Time: Aug 10 01:32:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 09 01:00:03 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	41398	20.000	ng/u1	0.00
20) Naphthalene-d8	10.748	136	189031	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.578	164	147280	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.315	188	310274	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	319920	20.000	ng/u1	0.00
88) Perylene-d12	24.641	264	368253	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.422	96	5921	6.167	ng/uL	0.00
4) Pyridine-d5	3.822	84	89687	29.694	ng/u1	0.00
7) Phenol-d5	7.106	99	132333	33.491	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.282	67	77065	31.684	ng/u1	0.00
11) 2-Chlorophenol-d4	7.488	132	91512	32.325	ng/u1	0.00
15) 4-Methylphenol-d8	8.651	113	107035	32.521	ng/u1	0.00
21) Nitrobenzene-d5	9.103	128	46016	39.478	ng/u1	0.00
24) 2-Nitrophenol-d4	9.826	143	53632	48.913	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.366	165	112631	35.367	ng/u1	0.00
31) 4-Chloroaniline-d4	10.877	131	155701	33.708	ng/u1	0.00
46) Dimethylphthalate-d6	13.985	166	364933	32.115	ng/u1	0.00
49) Acenaphthylene-d8	14.267	160	417910	32.437	ng/u1	0.00
54) 4-Nitrophenol-d4	14.754	143	53837	33.842	ng/u1	0.00
60) Fluorene-d10	15.565	176	335326	32.876	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.671	200	40463	44.034	ng/u1	0.00
73) Anthracene-d10	17.415	188	523330	34.970	ng/u1	0.00
81) Pyrene-d10	19.695	212	651790	34.277	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.435	264	675126	34.295	ng/u1	0.01

Target Compounds	Qvalue

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

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