

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070824\
 Data File : BG061951.D
 Acq On : 8 Jul 2024 21:34
 Operator : MA/JU
 Sample : PB161746BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK746

Quant Time: Jul 08 23:22:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 03 02:38:46 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.095	152	39624	20.000	ng/u1	0.00
20) Naphthalene-d8	10.903	136	204803	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.716	164	153864	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.454	188	358926	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.714	240	326023	20.000	ng/u1	-0.01
88) Perylene-d12	24.951	264	393917	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.482	96	6578	6.238	ng/uL	0.00
4) Pyridine-d5	3.899	84	86164	26.576	ng/u1	0.00
7) Phenol-d5	7.243	99	133082	30.690	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.413	67	79535	28.964	ng/u1	0.00
11) 2-Chlorophenol-d4	7.624	132	95188	32.002	ng/u1	0.00
15) 4-Methylphenol-d8	8.794	113	106970	31.331	ng/u1	0.00
21) Nitrobenzene-d5	9.258	128	52893	33.958	ng/u1	0.00
24) 2-Nitrophenol-d4	9.981	143	65505	36.826	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.521	165	120137	35.016	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.038	131	134022	26.906	ng/u1	0.00
46) Dimethylphthalate-d6	14.117	166	421847	33.348	ng/u1	0.00
49) Acenaphthylene-d8	14.411	160	461003	34.853	ng/u1	0.00
54) 4-Nitrophenol-d4	14.910	143	64924	32.139	ng/u1	0.00
60) Fluorene-d10	15.703	176	370908	34.830	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.815	200	65931	34.976	ng/u1	-0.01
73) Anthracene-d10	17.554	188	571487	34.103	ng/u1	0.00
81) Pyrene-d10	19.834	212	665293	34.693	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.734	264	709884	36.337	ng/u1	0.00

Target Compounds Qvalue

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

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