

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070824\
 Data File : BG061966.D
 Acq On : 9 Jul 2024 9:03
 Operator : MA/JU
 Sample : PB161730BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK730

Quant Time: Jul 10 00:25:10 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	8.093	152	28904	20.000	ng/u1	0.00
20) Naphthalene-d8	10.902	136	150198	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.715	164	121503	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.459	188	292823	20.000	ng/u1	0.00
79) Chrysene-d12	21.719	240	275664	20.000	ng/u1	# 0.00
88) Perylene-d12	24.962	264	336297	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.481	96	5585	7.261	ng/uL	0.00
4) Pyridine-d5	3.898	84	60769	25.695	ng/u1	0.00
7) Phenol-d5	7.242	99	89785	28.385	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.412	67	52068	25.993	ng/u1	0.00
11) 2-Chlorophenol-d4	7.623	132	63409	29.224	ng/u1	0.00
15) 4-Methylphenol-d8	8.793	113	69930	28.079	ng/u1	0.00
21) Nitrobenzene-d5	9.257	128	35505	31.082	ng/u1	0.00
24) 2-Nitrophenol-d4	9.979	143	42494	32.575	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.520	165	79029	31.409	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.037	131	87118	23.848	ng/u1	0.00
46) Dimethylphthalate-d6	14.116	166	286972	28.728	ng/u1	0.00
49) Acenaphthylene-d8	14.415	160	305947	29.291	ng/u1	0.00
54) 4-Nitrophenol-d4	14.909	143	46515	29.159	ng/u1	0.00
60) Fluorene-d10	15.708	176	255955	30.437	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.820	200	45730	29.736	ng/u1	0.00
73) Anthracene-d10	17.559	188	423321	30.964	ng/u1	0.00
81) Pyrene-d10	19.839	212	491724	30.327	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.739	264	518600	31.094	ng/u1	0.00

Target Compounds	Qvalue

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

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