

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070824\
 Data File : BG061977.D
 Acq On : 9 Jul 2024 17:19
 Operator : MA/JU
 Sample : PB161754BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK754

Quant Time: Jul 10 00:40:23 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	35801	20.000	ng/u1	0.00
20) Naphthalene-d8	10.909	136	183509	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.716	164	142276	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.454	188	324847	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.714	240	288841	20.000	ng/u1	-0.01
88) Perylene-d12	24.951	264	350715	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.488	96	5339	5.604	ng/uL	0.00
4) Pyridine-d5	3.905	84	73359	25.043	ng/u1	0.00
7) Phenol-d5	7.243	99	107336	27.396	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.419	67	65540	26.416	ng/u1	0.00
11) 2-Chlorophenol-d4	7.625	132	76116	28.322	ng/u1	0.00
15) 4-Methylphenol-d8	8.794	113	87127	28.244	ng/u1	0.00
21) Nitrobenzene-d5	9.258	128	42081	30.152	ng/u1	0.00
24) 2-Nitrophenol-d4	9.981	143	51076	32.047	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.521	165	97363	31.671	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.038	131	138419	31.014	ng/u1	0.00
46) Dimethylphthalate-d6	14.117	166	351917	30.086	ng/u1	0.00
49) Acenaphthylene-d8	14.411	160	376200	30.758	ng/u1	0.00
54) 4-Nitrophenol-d4	14.904	143	52367	28.034	ng/u1	-0.01
60) Fluorene-d10	15.703	176	299276	30.393	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.815	200	54753	32.094	ng/u1	-0.01
73) Anthracene-d10	17.554	188	478552	31.553	ng/u1	0.00
81) Pyrene-d10	19.834	212	542902	31.955	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.728	264	568057	32.659	ng/u1	-0.01

Target Compounds Qvalue

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

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