

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071024\
 Data File : BG061992.D
 Acq On : 10 Jul 2024 10:25
 Operator : MA/JU
 Sample : PB161916BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK916

Quant Time: Jul 11 00:41:48 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	8.094	152	54777	20.000	ng/u1	0.00
20) Naphthalene-d8	10.908	136	282724	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.716	164	210727	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.453	188	484865	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.713	240	435066	20.000	ng/u1	-0.01
88) Perylene-d12	24.956	264	541237	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.476	96	5509	3.779	ng/uL	-0.01
4) Pyridine-d5	3.899	84	89783	20.032	ng/u1	0.00
7) Phenol-d5	7.248	99	132527	22.108	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.412	67	81176	21.384	ng/u1	0.00
11) 2-Chlorophenol-d4	7.624	132	94388	22.954	ng/u1	0.00
15) 4-Methylphenol-d8	8.793	113	104475	22.135	ng/u1	0.00
21) Nitrobenzene-d5	9.263	128	49812	23.166	ng/u1	0.00
24) 2-Nitrophenol-d4	9.980	143	59590	24.268	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.520	165	115545	24.396	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.037	131	130394	18.963	ng/u1	0.00
46) Dimethylphthalate-d6	14.116	166	399182	23.041	ng/u1	0.00
49) Acenaphthylene-d8	14.410	160	432779	23.890	ng/u1	0.00
54) 4-Nitrophenol-d4	14.909	143	58997	21.324	ng/u1	0.00
60) Fluorene-d10	15.703	176	352113	24.143	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.814	200	61945	24.326	ng/u1	-0.01
73) Anthracene-d10	17.553	188	556935	24.602	ng/u1	0.00
81) Pyrene-d10	19.833	212	624545	24.406	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.727	264	668540	24.906	ng/u1	-0.01

Target Compounds Qvalue

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

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