

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071023\
 Data File : BG058138.D
 Acq On : 10 Jul 2023 12:06
 Operator : CG/JU
 Sample : PB153911BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK911

Quant Time: Jul 11 00:44:38 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 10 16:52:40 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.972	152	55110	20.000	ng/ul	0.00
20) Naphthalene-d8	10.775	136	251728	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.600	164	183105	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.344	188	452948	20.000	ng/ul	0.00
79) Chrysene-d12	21.603	240	418688	20.000	ng/ul	0.00
88) Perylene-d12	24.794	264	508045	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.401	96	6838	4.345	ng/uL	0.00
4) Pyridine-d5	3.812	84	109492	22.916	ng/ul	0.00
7) Phenol-d5	7.132	99	122440	21.755	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.297	67	76985	22.698	ng/ul	0.00
11) 2-Chlorophenol-d4	7.502	132	86365	22.249	ng/ul	0.00
15) 4-Methylphenol-d8	8.671	113	93905	22.112	ng/ul	0.00
21) Nitrobenzene-d5	9.136	128	45652	22.251	ng/ul	0.00
24) 2-Nitrophenol-d4	9.852	143	50654	23.491	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	88739	21.452	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	132767	21.170	ng/ul	0.00
46) Dimethylphthalate-d6	14.000	166	324935	22.491	ng/ul	0.00
49) Acenaphthylene-d8	14.294	160	357786	22.695	ng/ul	0.00
54) 4-Nitrophenol-d4	14.800	143	48183	18.877	ng/ul	0.00
60) Fluorene-d10	15.593	176	284377	23.178	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	47083	19.143	ng/ul	0.00
73) Anthracene-d10	17.443	188	480597	22.866	ng/ul	0.00
81) Pyrene-d10	19.729	212	577294	21.893	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.576	264	596476	23.044	ng/ul	0.00

Target Compounds Qvalue

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

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