

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
 Data File : BG059402.D
 Acq On : 9 Oct 2023 18:34
 Operator : MA/JU
 Sample : PB156170BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK170

Quant Time: Oct 09 22:46:59 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 07 22:45:22 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.226	152	44307	20.000	ng/ul	0.00
20) Naphthalene-d8	11.064	136	213237	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.859	164	148183	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.609	188	349962	20.000	ng/ul	0.00
79) Chrysene-d12	21.910	240	305321	20.000	ng/ul	0.00
88) Perylene-d12	25.388	264	371197	20.000	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.525	96	7889	6.348	ng/uL	0.00
4) Pyridine-d5	3.960	84	120867	33.792	ng/ul	0.00
7) Phenol-d5	7.356	99	160493	32.805	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.544	67	92144	33.323	ng/ul	0.00
11) 2-Chlorophenol-d4	7.744	132	115468	33.747	ng/ul	0.00
15) 4-Methylphenol-d8	8.919	113	121578	31.896	ng/ul	0.00
21) Nitrobenzene-d5	9.424	128	63631	35.054	ng/ul	0.00
24) 2-Nitrophenol-d4	10.141	143	69489	35.369	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.670	165	119498	33.698	ng/ul	0.00
31) 4-Chloroaniline-d4	11.216	131	187709	33.324	ng/ul	0.00
46) Dimethylphthalate-d6	14.254	166	422196	34.777	ng/ul	0.00
49) Acenaphthylene-d8	14.565	160	492892	34.117	ng/ul	0.00
54) 4-Nitrophenol-d4	15.053	143	66102	29.056	ng/ul	0.00
60) Fluorene-d10	15.846	176	386841	33.383	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.958	200	63838	27.628	ng/ul	0.00
73) Anthracene-d10	17.709	188	604092	35.188	ng/ul	0.00
81) Pyrene-d10	19.982	212	684502	37.008	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.147	264	705578	36.122	ng/ul	0.00

Target Compounds Qvalue

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

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