

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101323\
 Data File : BG059441.D
 Acq On : 13 Oct 2023 11:53
 Operator : MA/JU
 Sample : PB156220BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK220

Quant Time: Oct 14 00:50:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 12 17:05:49 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.220	152	37764	20.000	ng/ul	0.00
20) Naphthalene-d8	11.058	136	167022	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.859	164	104034	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.609	188	241532	20.000	ng/ul	0.00
79) Chrysene-d12	21.910	240	248666	20.000	ng/ul	0.00
88) Perylene-d12	25.388	264	292361	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.508	96	6261	5.199	ng/uL	0.00
4) Pyridine-d5	3.942	84	87746	26.623	ng/ul	0.00
7) Phenol-d5	7.344	99	110680	26.768	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.538	67	68350	29.004	ng/ul	0.00
11) 2-Chlorophenol-d4	7.732	132	86173	29.289	ng/ul	0.00
15) 4-Methylphenol-d8	8.907	113	83132	26.368	ng/ul	0.00
21) Nitrobenzene-d5	9.418	128	42516	30.824	ng/ul	0.00
24) 2-Nitrophenol-d4	10.135	143	41766	31.018	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.664	165	76935	27.495	ng/ul	0.00
31) 4-Chloroaniline-d4	11.204	131	122559	27.600	ng/ul	0.00
46) Dimethylphthalate-d6	14.254	166	287854	32.658	ng/ul	0.00
49) Acenaphthylene-d8	14.559	160	319056	30.021	ng/ul	0.00
54) 4-Nitrophenol-d4	15.047	143	39594	26.006	ng/ul	0.00
60) Fluorene-d10	15.846	176	248611	29.965	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.958	200	39727	26.918	ng/ul	0.00
73) Anthracene-d10	17.709	188	402523	32.415	ng/ul	0.00
81) Pyrene-d10	19.982	212	484113	32.053	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.147	264	535044	33.428	ng/ul	0.00

Target Compounds Qvalue

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

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