

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092823\
 Data File : BG059227.D
 Acq On : 28 Sep 2023 10:12
 Operator : MA/JU
 Sample : PB155787BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK787

Quant Time: Sep 29 00:41:08 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 28 04:33:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.233	152	50526	20.000	ng/ul	0.00
20) Naphthalene-d8	11.070	136	257833	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.866	164	165454	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.616	188	378407	20.000	ng/ul	0.00
79) Chrysene-d12	21.917	240	359269	20.000	ng/ul	0.00
88) Perylene-d12	25.401	264	420746	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.526	96	9170	6.850	ng/uL	0.00
4) Pyridine-d5	3.961	84	128646	32.934	ng/ul	0.00
7) Phenol-d5	7.357	99	154129	30.475	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.551	67	104731	32.416	ng/ul	0.00
11) 2-Chlorophenol-d4	7.745	132	110780	32.585	ng/ul	0.00
15) 4-Methylphenol-d8	8.926	113	117195	29.977	ng/ul	0.00
21) Nitrobenzene-d5	9.431	128	57220	31.449	ng/ul	0.00
24) 2-Nitrophenol-d4	10.148	143	61479	34.421	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.677	165	110259	30.050	ng/ul	0.00
31) 4-Chloroaniline-d4	11.217	131	179125	30.174	ng/ul	0.00
46) Dimethylphthalate-d6	14.261	166	399188	33.370	ng/ul	0.00
49) Acenaphthylene-d8	14.566	160	452451	30.134	ng/ul	0.00
54) 4-Nitrophenol-d4	15.054	143	61029	29.215	ng/ul	0.00
60) Fluorene-d10	15.853	176	349121	30.664	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.965	200	56709	29.155	ng/ul	0.00
73) Anthracene-d10	17.716	188	542054	31.438	ng/ul	0.00
81) Pyrene-d10	19.989	212	645368	32.077	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.160	264	653635	32.328	ng/ul	0.00

Target Compounds Qvalue

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

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