

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090823\
 Data File : BG058925.D
 Acq On : 8 Sep 2023 20:10
 Operator : MA/JU
 Sample : PB155275BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Sep 09 03:45:22 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 09 00:11:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.335	152	56922	20.000	ng/ul	0.00
20) Naphthalene-d8	11.179	136	251427	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.957	164	236292	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.707	188	675592	20.000	ng/ul	0.00
79) Chrysene-d12	22.031	240	569046	20.000	ng/ul	0.00
88) Perylene-d12	25.609	264	669626	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.641	96	6948	5.855	ng/uL	0.00
4) Pyridine-d5	4.070	84	111625	27.394	ng/ul	0.00
7) Phenol-d5	7.436	99	152183	26.956	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.648	67	88807	27.806	ng/ul	0.00
11) 2-Chlorophenol-d4	7.842	132	95203	28.270	ng/ul	0.00
15) 4-Methylphenol-d8	9.011	113	114357	26.135	ng/ul	0.00
21) Nitrobenzene-d5	9.522	128	48160	32.519	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.251	143	54869	33.821	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.768	165	133240	28.308	ng/ul	0.00
31) 4-Chloroaniline-d4	11.314	131	168470	27.903	ng/ul	0.00
46) Dimethylphthalate-d6	14.358	166	571680	31.342	ng/ul	0.00
49) Acenaphthylene-d8	14.657	160	563238	28.568	ng/ul	0.00
54) 4-Nitrophenol-d4	15.115	143	73115	30.339	ng/ul	0.00
60) Fluorene-d10	15.944	176	502296	29.218	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.044	200	87353	28.400	ng/ul	0.00
73) Anthracene-d10	17.806	188	917679	29.861	ng/ul	0.00
81) Pyrene-d10	20.074	212	1010251	30.612	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.356	264	1017962	30.860	ng/ul	0.00

Target Compounds	Qvalue

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

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