

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092823\
 Data File : BG059226.D
 Acq On : 28 Sep 2023 9:31
 Operator : MA/JU
 Sample : PB155814BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK814

Quant Time: Sep 29 00:55:17 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	8.237	152	41444	20.000	ng/u1	0.00
20) Naphthalene-d8	11.075	136	207239	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.870	164	139574	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.620	188	310457	20.000	ng/u1	0.00
79) Chrysene-d12	21.921	240	292051	20.000	ng/u1	0.00
88) Perylene-d12	25.405	264	347701	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.531	96	8929	8.131	ng/uL	0.00
4) Pyridine-d5	3.960	84	128637	40.149	ng/u1	0.00
7) Phenol-d5	7.356	99	161815	39.006	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.555	67	107619	40.609	ng/u1	0.00
11) 2-Chlorophenol-d4	7.749	132	114336	41.000	ng/u1	0.00
15) 4-Methylphenol-d8	8.924	113	124074	38.691	ng/u1	0.00
21) Nitrobenzene-d5	9.430	128	59612	40.763	ng/u1	0.00
24) 2-Nitrophenol-d4	10.147	143	64964	45.251	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.675	165	114973	38.985	ng/u1	0.00
31) 4-Chloroaniline-d4	11.222	131	186046	38.991	ng/u1	0.00
46) Dimethylphthalate-d6	14.265	166	439659	43.569	ng/u1	0.00
49) Acenaphthylene-d8	14.571	160	488100	38.536	ng/u1	0.00
54) 4-Nitrophenol-d4	15.058	143	66189	37.561	ng/u1	0.00
60) Fluorene-d10	15.858	176	375553	39.102	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.963	200	62766	39.332	ng/u1	0.00
73) Anthracene-d10	17.720	188	599249	42.362	ng/u1	0.00
81) Pyrene-d10	19.994	212	717051	43.843	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.164	264	734636	43.967	ng/u1	0.00

Target Compounds	Qvalue

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

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