

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092823\
 Data File : BG059241.D
 Acq On : 29 Sep 2023 5:52
 Operator : MA/JU
 Sample : PB155826BL
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK826

Quant Time: Sep 29 06:29:28 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	8.226	152	39237	20.000	ng/u1	0.00
20) Naphthalene-d8	11.076	136	181090	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.865	164	118881	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.615	188	271087	20.000	ng/u1	0.00
79) Chrysene-d12	21.916	240	257292	20.000	ng/u1	0.00
88) Perylene-d12	25.400	264	297707	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.526	96	7998	7.693	ng/uL	0.00
4) Pyridine-d5	3.960	84	109209	36.002	ng/u1	0.00
7) Phenol-d5	7.356	99	146151	37.212	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.550	67	93181	37.139	ng/u1	0.00
11) 2-Chlorophenol-d4	7.744	132	107226	40.614	ng/u1	0.00
15) 4-Methylphenol-d8	8.919	113	110231	36.308	ng/u1	-0.01
21) Nitrobenzene-d5	9.430	128	52743	41.274	ng/u1	0.00
24) 2-Nitrophenol-d4	10.147	143	57316	45.689	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.676	165	104897	40.704	ng/u1	0.00
31) 4-Chloroaniline-d4	11.217	131	159438	38.239	ng/u1	0.00
46) Dimethylphthalate-d6	14.260	166	362586	42.185	ng/u1	0.00
49) Acenaphthylene-d8	14.566	160	408040	37.822	ng/u1	0.00
54) 4-Nitrophenol-d4	15.053	143	56892	37.904	ng/u1	0.00
60) Fluorene-d10	15.852	176	322622	39.438	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.958	200	43397	31.144	ng/u1	0.00
73) Anthracene-d10	17.715	188	523966	42.419	ng/u1	0.00
81) Pyrene-d10	19.989	212	609949	42.332	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.159	264	625688	43.735	ng/u1	0.00

Target Compounds Qvalue

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

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