

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100723\
 Data File : BG059347.D
 Acq On : 7 Oct 2023 18:12
 Operator : MA/JU
 Sample : PB155975BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK975

Quant Time: Oct 07 23:45:13 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 07 22:45:22 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.231	152	47306	20.000	ng/u1	0.00
20) Naphthalene-d8	11.069	136	233518	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.865	164	165785	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.615	188	418523	20.000	ng/u1	0.00
79) Chrysene-d12	21.915	240	395047	20.000	ng/u1	0.00
88) Perylene-d12	25.394	264	463349	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.525	96	9403	7.087	ng/uL	0.00
4) Pyridine-d5	3.960	84	131832	34.521	ng/u1	0.00
7) Phenol-d5	7.356	99	163361	31.275	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.550	67	98044	33.209	ng/u1	0.00
11) 2-Chlorophenol-d4	7.744	132	122245	33.463	ng/u1	0.00
15) 4-Methylphenol-d8	8.919	113	123526	30.352	ng/u1	0.00
21) Nitrobenzene-d5	9.430	128	66079	33.241	ng/u1	0.00
24) 2-Nitrophenol-d4	10.147	143	73548	34.184	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.676	165	123641	31.838	ng/u1	0.00
31) 4-Chloroaniline-d4	11.216	131	201467	32.660	ng/u1	0.00
46) Dimethylphthalate-d6	14.260	166	481511	35.452	ng/u1	0.00
49) Acenaphthylene-d8	14.565	160	541596	33.508	ng/u1	0.00
54) 4-Nitrophenol-d4	15.053	143	82684	32.486	ng/u1	0.00
60) Fluorene-d10	15.852	176	438213	33.801	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.964	200	83799	30.326	ng/u1	0.00
73) Anthracene-d10	17.714	188	717585	34.952	ng/u1	0.00
81) Pyrene-d10	19.988	212	820836	34.299	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.153	264	840553	34.474	ng/u1	0.00

Target Compounds Qvalue

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

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