

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
 Data File : BG059393.D
 Acq On : 9 Oct 2023 11:35
 Operator : MA/JU
 Sample : PB155981BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK981

Quant Time: Oct 09 23:14:17 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	8.226	152	42809	20.000	ng/u1	0.00
20) Naphthalene-d8	11.070	136	214604	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.859	164	152487	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.609	188	356634	20.000	ng/u1	0.00
79) Chrysene-d12	21.910	240	315711	20.000	ng/u1	0.00
88) Perylene-d12	25.400	264	370426	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.525	96	10373	8.639	ng/uL	0.00
4) Pyridine-d5	3.960	84	140845	40.755	ng/u1	0.00
7) Phenol-d5	7.350	99	183162	38.749	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.544	67	107700	40.311	ng/u1	0.00
11) 2-Chlorophenol-d4	7.744	132	135915	41.113	ng/u1	0.00
15) 4-Methylphenol-d8	8.919	113	139027	37.750	ng/u1	0.00
21) Nitrobenzene-d5	9.424	128	73972	40.491	ng/u1	0.00
24) 2-Nitrophenol-d4	10.141	143	82358	41.652	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.670	165	140996	39.507	ng/u1	0.00
31) 4-Chloroaniline-d4	11.216	131	218316	38.510	ng/u1	0.00
46) Dimethylphthalate-d6	14.260	166	508162	40.677	ng/u1	0.00
49) Acenaphthylene-d8	14.565	160	582062	39.152	ng/u1	0.00
54) 4-Nitrophenol-d4	15.053	143	78598	33.573	ng/u1	0.00
60) Fluorene-d10	15.852	176	463279	38.851	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.964	200	77814	33.047	ng/u1	0.00
73) Anthracene-d10	17.709	188	725853	41.490	ng/u1	0.00
81) Pyrene-d10	19.988	212	791047	41.361	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.153	264	809879	41.548	ng/u1	0.00

Target Compounds Qvalue

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

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