

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092823\
 Data File : BG059231.D
 Acq On : 28 Sep 2023 12:54
 Operator : MA/JU
 Sample : PB155590TB 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SLEB590

Quant Time: Sep 29 00:41:42 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	8.229	152	41817	20.000	ng/ul	0.00
20) Naphthalene-d8	11.073	136	227141	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.868	164	149835	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.618	188	344215	20.000	ng/ul	0.00
79) Chrysene-d12	21.919	240	302538	20.000	ng/ul	0.00
88) Perylene-d12	25.403	264	355444	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.529	96	7358	6.641	ng/uL	0.00
4) Pyridine-d5	3.963	84	111450	34.474	ng/ul	0.00
7) Phenol-d5	7.354	99	140416	33.546	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.553	67	94467	35.329	ng/ul	0.00
11) 2-Chlorophenol-d4	7.747	132	101043	35.910	ng/ul	0.00
15) 4-Methylphenol-d8	8.922	113	108475	33.525	ng/ul	0.00
21) Nitrobenzene-d5	9.428	128	54445	33.968	ng/ul	0.00
24) 2-Nitrophenol-d4	10.144	143	56961	36.200	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.673	165	99717	30.849	ng/ul	0.00
31) 4-Chloroaniline-d4	11.220	131	164163	31.390	ng/ul	0.00
46) Dimethylphthalate-d6	14.263	166	367197	33.896	ng/ul	0.00
49) Acenaphthylene-d8	14.569	160	427493	31.439	ng/ul	0.00
54) 4-Nitrophenol-d4	15.056	143	56002	29.603	ng/ul	0.00
60) Fluorene-d10	15.855	176	329217	31.930	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.961	200	52146	29.472	ng/ul	0.00
73) Anthracene-d10	17.718	188	513545	32.743	ng/ul	0.00
81) Pyrene-d10	19.992	212	593862	35.052	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.156	264	584372	34.212	ng/ul	0.00

Target Compounds Qvalue

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

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