

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091323\
 Data File : BG059051.D
 Acq On : 13 Sep 2023 22:51
 Operator : MA/JU
 Sample : PB155393TB 10X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SLEB393

Quant Time: Sep 14 01:07:44 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 12 23:54:19 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.302	152	80883	20.000	ng/u1	0.00
20) Naphthalene-d8	11.146	136	371418	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.930	164	306312	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.679	188	848563	20.000	ng/u1	0.00
79) Chrysene-d12	22.004	240	768017	20.000	ng/u1	0.00
88) Perylene-d12	25.552	264	852203	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.613	96	11711	6.154	ng/uL	-0.02
4) Pyridine-d5	4.042	84	146817	24.369	ng/u1	-0.02
7) Phenol-d5	7.415	99	254692	31.226	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.620	67	144636	30.242	ng/u1	0.00
11) 2-Chlorophenol-d4	7.814	132	168101	32.135	ng/u1	0.00
15) 4-Methylphenol-d8	8.984	113	204387	32.435	ng/u1	0.00
21) Nitrobenzene-d5	9.501	128	91882	34.105	ng/u1	0.00
24) 2-Nitrophenol-d4	10.212	143	118442	40.066	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.734	165	209936	33.415	ng/u1	-0.02
31) 4-Chloroaniline-d4	11.287	131	185132	21.912	ng/u1	0.00
46) Dimethylphthalate-d6	14.324	166	834902	35.139	ng/u1	0.00
49) Acenaphthylene-d8	14.630	160	857424	32.406	ng/u1	0.00
54) 4-Nitrophenol-d4	15.100	143	143469	36.351	ng/u1	0.00
60) Fluorene-d10	15.917	176	759696	36.028	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	16.022	200	160726	38.587	ng/u1	0.00
73) Anthracene-d10	17.779	188	1301310	34.545	ng/u1	0.00
81) Pyrene-d10	20.053	212	1499445	36.856	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.311	264	1443887	34.563	ng/u1	0.00

Target Compounds	Qvalue

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

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