

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101623\
 Data File : BG059460.D
 Acq On : 16 Oct 2023 17:09
 Operator : MA/JU
 Sample : PB156298BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK298

Quant Time: Oct 17 03:34:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 12 17:05:49 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.217	152	31936	20.000	ng/ul	0.00
20) Naphthalene-d8	11.061	136	141321	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.856	164	87925	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.606	188	217337	20.000	ng/ul	0.00
79) Chrysene-d12	21.907	240	235170	20.000	ng/ul	0.00
88) Perylene-d12	25.385	264	279023	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.517	96	7211	7.081	ng/uL	0.00
4) Pyridine-d5	3.945	84	132532	47.550	ng/ul	0.00
7) Phenol-d5	7.347	99	108544	31.042	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.535	67	64634	32.433	ng/ul	0.00
11) 2-Chlorophenol-d4	7.735	132	85086	34.197	ng/ul	0.00
15) 4-Methylphenol-d8	8.910	113	80128	30.053	ng/ul	0.00
21) Nitrobenzene-d5	9.415	128	57079	48.907	ng/ul	0.00
24) 2-Nitrophenol-d4	10.132	143	54252	47.619	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.661	165	73609	31.091	ng/ul	0.00
31) 4-Chloroaniline-d4	11.208	131	120900	32.178	ng/ul	0.00
46) Dimethylphthalate-d6	14.251	166	268293	36.015	ng/ul	0.00
49) Acenaphthylene-d8	14.557	160	304892	33.944	ng/ul	0.00
54) 4-Nitrophenol-d4	15.044	143	41474	32.232	ng/ul	0.00
60) Fluorene-d10	15.843	176	236242	33.691	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.955	200	42747	32.189	ng/ul	0.00
73) Anthracene-d10	17.706	188	386771	34.614	ng/ul	0.00
81) Pyrene-d10	19.980	212	504924	35.349	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.144	264	559951	36.656	ng/ul	0.00

Target Compounds Qvalue

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

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