

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101323\
 Data File : BG059445.D
 Acq On : 13 Oct 2023 16:36
 Operator : MA/JU
 Sample : PB156222BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK222

Quant Time: Oct 14 00:52:29 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	40391	20.000	ng/ul	0.00
20) Naphthalene-d8	11.061	136	181338	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.856	164	113879	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.606	188	269661	20.000	ng/ul	0.00
79) Chrysene-d12	21.907	240	286163	20.000	ng/ul	0.00
88) Perylene-d12	25.385	264	339419	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.511	96	7282	5.654	ng/uL	0.00
4) Pyridine-d5	3.946	84	89440	25.372	ng/ul	0.00
7) Phenol-d5	7.342	99	113931	25.762	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.536	67	68431	27.150	ng/ul	0.00
11) 2-Chlorophenol-d4	7.735	132	82243	26.135	ng/ul	0.00
15) 4-Methylphenol-d8	8.910	113	81472	24.161	ng/ul	0.00
21) Nitrobenzene-d5	9.416	128	41039	27.404	ng/ul	0.00
24) 2-Nitrophenol-d4	10.138	143	43952	30.065	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.661	165	77340	25.458	ng/ul	0.00
31) 4-Chloroaniline-d4	11.208	131	124232	25.768	ng/ul	0.00
46) Dimethylphthalate-d6	14.251	166	274225	28.422	ng/ul	0.00
49) Acenaphthylene-d8	14.557	160	308058	26.480	ng/ul	0.00
54) 4-Nitrophenol-d4	15.044	143	38837	23.304	ng/ul	0.00
60) Fluorene-d10	15.844	176	245320	27.012	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.955	200	40040	24.301	ng/ul	0.00
73) Anthracene-d10	17.706	188	387484	27.949	ng/ul	0.00
81) Pyrene-d10	19.980	212	475809	27.375	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.150	264	505697	27.214	ng/ul	0.00

Target Compounds Qvalue

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

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