

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100323\
 Data File : BG059267.D
 Acq On : 3 Oct 2023 16:33
 Operator : MA/JU
 Sample : 04480-17
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BBJN2

Quant Time: Oct 03 23:33:49 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	8.235	152	44313	20.000	ng/ul	0.00
20) Naphthalene-d8	11.073	136	219817	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.862	164	143395	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.612	188	322423	20.000	ng/ul	0.00
79) Chrysene-d12	21.913	240	291723	20.000	ng/ul	0.00
88) Perylene-d12	25.391	264	355595	20.000	ng/ul	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.534	96	4917	4.188	ng/uL	0.00
4) Pyridine-d5	3.963	84	21355	6.234	ng/ul	0.00
7) Phenol-d5	7.359	99	26805	6.043	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.553	67	73157	25.818	ng/ul	0.00
11) 2-Chlorophenol-d4	7.747	132	65680	22.028	ng/ul	0.00
15) 4-Methylphenol-d8	8.922	113	41941	12.232	ng/ul	0.00
21) Nitrobenzene-d5	9.427	128	43015	27.731	ng/ul	0.00
24) 2-Nitrophenol-d4	10.150	143	43442	28.528	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.679	165	79576	25.439	ng/ul	0.00
31) 4-Chloroaniline-d4	11.220	131	78915	15.592	ng/ul	0.00
46) Dimethylphthalate-d6	14.263	166	313273	30.217	ng/ul	0.00
49) Acenaphthylene-d8	14.569	160	355071	27.286	ng/ul	0.00
54) 4-Nitrophenol-d4	15.056	143	10803	5.967	ng/ul	0.00
60) Fluorene-d10	15.849	176	285949	28.979	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.961	200	42727	25.781	ng/ul	0.00
73) Anthracene-d10	17.712	188	462488	31.481	ng/ul	0.00
81) Pyrene-d10	19.986	212	543333	33.258	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.144	264	539433	31.568	ng/ul	-0.02

Target Compounds	Qvalue

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

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