

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100223\
 Data File : BG059252.D
 Acq On : 2 Oct 2023 16:57
 Operator : MA/JU
 Sample : 04480-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BBJ47

Quant Time: Oct 02 23:55:43 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	53524	20.000	ng/u1	0.00
20) Naphthalene-d8	11.073	136	255238	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.863	164	166073	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.613	188	366511	20.000	ng/u1	0.00
79) Chrysene-d12	21.907	240	327546	20.000	ng/u1	-0.01
88) Perylene-d12	25.392	264	386942	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.529	96	5610	3.956	ng/uL	0.00
4) Pyridine-d5	3.964	84	25846	6.246	ng/u1	0.00
7) Phenol-d5	7.354	99	27059	5.051	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.554	67	70670	20.648	ng/u1	0.00
11) 2-Chlorophenol-d4	7.748	132	61222	16.999	ng/u1	0.00
15) 4-Methylphenol-d8	8.923	113	45537	10.995	ng/u1	0.00
21) Nitrobenzene-d5	9.428	128	42214	23.438	ng/u1	0.00
24) 2-Nitrophenol-d4	10.145	143	44864	25.374	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.674	165	74373	20.476	ng/u1	0.00
31) 4-Chloroaniline-d4	11.214	131	74158	12.619	ng/u1	0.00
46) Dimethylphthalate-d6	14.258	166	285929	23.813	ng/u1	0.00
49) Acenaphthylene-d8	14.569	160	330968	21.961	ng/u1	0.00
54) 4-Nitrophenol-d4	15.057	143	8569	4.087	ng/u1	0.00
60) Fluorene-d10	15.850	176	260870	22.827	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.961	200	32633	17.322	ng/u1	0.00
73) Anthracene-d10	17.712	188	390318	23.372	ng/u1	0.00
81) Pyrene-d10	19.986	212	449267	24.493	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.139	264	433580	23.318	ng/u1	-0.03

Target Compounds Qvalue

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

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