

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
 Data File : BG050958.D
 Acq On : 11 Nov 2021 1:51
 Operator : CG/JU
 Sample : M4445-21
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG3A4

Quant Time: Nov 11 03:56:41 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 02 14:49:05 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.224	152	40060	20.000	ng/u1	-0.01
20) Naphthalene-d8	11.050	136	172947	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.845	164	118042	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.595	188	261334	20.000	ng/u1	-0.01
79) Chrysene-d12	21.890	240	228412	20.000	ng/u1	-0.02
88) Perylene-d12	25.286	264	222826	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.588	96	4959	3.995	ng/uL	0.00
4) Pyridine-d5	4.017	84	20799	5.601	ng/u1	0.00
7) Phenol-d5	7.372	99	20836	4.875	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	67211	24.346	ng/u1	0.00
11) 2-Chlorophenol-d4	7.754	132	55809	18.843	ng/u1	-0.01
15) 4-Methylphenol-d8	8.923	113	38240	11.366	ng/u1	-0.01
21) Nitrobenzene-d5	9.399	128	37638	25.609	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.127	143	38395	23.495	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.662	165	60569	22.003	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.185	131	56619	13.582	ng/u1	0.00
46) Dimethylphthalate-d6	14.240	166	259380	28.721	ng/u1	-0.01
49) Acenaphthylene-d8	14.546	160	298996	26.574	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.051	143	6824	4.167	ng/u1	0.00
60) Fluorene-d10	15.838	176	230670	28.833	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.950	200	37838	23.881	ng/u1	-0.01
73) Anthracene-d10	17.689	188	410431	33.218	ng/u1	-0.02
81) Pyrene-d10	19.969	212	484024	32.809	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.051	264	404287	32.820	ng/u1	-0.02

Target Compounds	Qvalue

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

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