

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
 Data File : BG050883.D
 Acq On : 8 Nov 2021 16:53
 Operator : CG/JU
 Sample : M4492-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG1R8

Quant Time: Nov 08 16:42:39 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	8.230	152	31949	20.000	ng/u1	0.00
20) Naphthalene-d8	11.056	136	143195	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.851	164	98039	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.595	188	217431	20.000	ng/u1	-0.01
79) Chrysene-d12	21.890	240	193495	20.000	ng/u1	-0.02
88) Perylene-d12	25.286	264	187633	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.588	96	2571	2.597	ng/uL	0.00
4) Pyridine-d5	4.023	84	22700	7.665	ng/u1	0.00
7) Phenol-d5	7.372	99	24585	7.213	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	68303	31.022	ng/u1	0.00
11) 2-Chlorophenol-d4	7.760	132	57267	24.244	ng/u1	0.00
15) 4-Methylphenol-d8	8.929	113	41419	15.436	ng/u1	0.00
21) Nitrobenzene-d5	9.399	128	38134	31.338	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.128	143	38765	28.650	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.668	165	61106	26.810	ng/u1	0.00
31) 4-Chloroaniline-d4	11.185	131	93128	26.981	ng/u1	0.00
46) Dimethylphthalate-d6	14.246	166	260803	34.771	ng/u1	0.00
49) Acenaphthylene-d8	14.546	160	310409	33.217	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.039	143	8655	6.364	ng/u1	0.00
60) Fluorene-d10	15.839	176	236969	35.664	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.956	200	38654	29.323	ng/u1	0.00
73) Anthracene-d10	17.695	188	420834	40.938	ng/u1	-0.01
81) Pyrene-d10	19.969	212	483086	38.654	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.057	264	405595	39.102	ng/u1	-0.02

Target Compounds	Qvalue

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

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