

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
 Data File : BG050919.D
 Acq On : 9 Nov 2021 21:44
 Operator : CG/JU
 Sample : M4445-17
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG391

Quant Time: Nov 10 07:14:58 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	45210	20.000	ng/u1	0.00
20) Naphthalene-d8	11.050	136	200244	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.851	164	132601	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.595	188	293050	20.000	ng/u1	-0.01
79) Chrysene-d12	21.890	240	252810	20.000	ng/u1	-0.02
88) Perylene-d12	25.292	264	254200	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.588	96	2771	1.978	ng/uL	0.00
4) Pyridine-d5	4.023	84	16973	4.050	ng/u1	0.00
7) Phenol-d5	7.372	99	29131	6.040	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	87077	27.949	ng/u1	0.00
11) 2-Chlorophenol-d4	7.754	132	70578	21.115	ng/u1	-0.01
15) 4-Methylphenol-d8	8.923	113	51038	13.442	ng/u1	-0.01
21) Nitrobenzene-d5	9.399	128	48825	28.692	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.127	143	50348	26.609	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.668	165	81431	25.549	ng/u1	0.00
31) 4-Chloroaniline-d4	11.185	131	87712	18.172	ng/u1	0.00
46) Dimethylphthalate-d6	14.246	166	330140	32.543	ng/u1	0.00
49) Acenaphthylene-d8	14.546	160	391993	31.014	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.039	143	10756	5.848	ng/u1	0.00
60) Fluorene-d10	15.838	176	285273	31.743	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.956	200	51215	28.826	ng/u1	0.00
73) Anthracene-d10	17.695	188	471100	34.002	ng/u1	-0.01
81) Pyrene-d10	19.969	212	561107	34.363	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.057	264	489177	34.810	ng/u1	-0.02

Target Compounds Qvalue

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

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