

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
 Data File : BG050904.D
 Acq On : 9 Nov 2021 7:57
 Operator : CG/JU
 Sample : M4445-11
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG385

Quant Time: Nov 09 07:35:41 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	8.229	152	32911	20.000	ng/u1	0.00
20) Naphthalene-d8	11.055	136	144486	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.850	164	96751	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.594	188	210811	20.000	ng/u1	-0.01
79) Chrysene-d12	21.889	240	185214	20.000	ng/u1	-0.02
88) Perylene-d12	25.291	264	181089	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.593	96	5463	5.357	ng/uL	0.00
4) Pyridine-d5	4.016	84	27711	9.084	ng/u1	0.00
7) Phenol-d5	7.371	99	18198	5.183	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	58376	25.739	ng/u1	0.00
11) 2-Chlorophenol-d4	7.759	132	46088	18.941	ng/u1	0.00
15) 4-Methylphenol-d8	8.928	113	32397	11.721	ng/u1	0.00
21) Nitrobenzene-d5	9.398	128	31989	26.053	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.127	143	32988	24.162	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.667	165	51282	22.299	ng/u1	0.00
31) 4-Chloroaniline-d4	11.184	131	74841	21.489	ng/u1	0.00
46) Dimethylphthalate-d6	14.245	166	210993	28.505	ng/u1	0.00
49) Acenaphthylene-d8	14.545	160	255599	27.716	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.044	143	5687	4.237	ng/u1	0.00
60) Fluorene-d10	15.837	176	188086	28.684	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.955	200	27446	21.474	ng/u1	0.00
73) Anthracene-d10	17.694	188	321167	32.223	ng/u1	-0.01
81) Pyrene-d10	19.968	212	374452	31.302	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.050	264	308424	30.809	ng/u1	-0.02

Target Compounds Qvalue

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

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