

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
 Data File : BG050903.D
 Acq On : 9 Nov 2021 7:16
 Operator : CG/JU
 Sample : M4445-12
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG386

Quant Time: Nov 09 07:33:32 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	35483	20.000	ng/u1	0.00
20) Naphthalene-d8	11.055	136	157629	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.851	164	103633	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.595	188	231037	20.000	ng/u1	-0.01
79) Chrysene-d12	21.890	240	205975	20.000	ng/u1	-0.02
88) Perylene-d12	25.292	264	197482	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.588	96	5437	4.946	ng/uL	0.00
4) Pyridine-d5	4.022	84	21384	6.502	ng/u1	0.00
7) Phenol-d5	7.371	99	21341	5.638	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	62846	25.701	ng/u1	0.00
11) 2-Chlorophenol-d4	7.753	132	51332	19.567	ng/u1	-0.01
15) 4-Methylphenol-d8	8.922	113	37286	12.512	ng/u1	-0.01
21) Nitrobenzene-d5	9.398	128	34826	25.998	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.127	143	35783	24.024	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.667	165	55271	22.029	ng/u1	0.00
31) 4-Chloroaniline-d4	11.185	131	79083	20.814	ng/u1	0.00
46) Dimethylphthalate-d6	14.246	166	236680	29.852	ng/u1	0.00
49) Acenaphthylene-d8	14.545	160	276118	27.953	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.039	143	6865	4.775	ng/u1	0.00
60) Fluorene-d10	15.838	176	206388	29.385	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.955	200	33484	23.905	ng/u1	0.00
73) Anthracene-d10	17.695	188	363752	33.301	ng/u1	-0.01
81) Pyrene-d10	19.968	212	430323	32.346	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.051	264	355932	32.603	ng/u1	-0.02

Target Compounds Qvalue

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

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