

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
 Data File : BG050889.D
 Acq On : 8 Nov 2021 21:01
 Operator : CG/JU
 Sample : M4445-09
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG383

Quant Time: Nov 09 01:05:17 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	35749	20.000	ng/u1	0.00
20) Naphthalene-d8	11.056	136	163111	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.851	164	108247	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.595	188	242030	20.000	ng/u1	-0.01
79) Chrysene-d12	21.890	240	207667	20.000	ng/u1	#-0.02
88) Perylene-d12	25.286	264	202631	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.594	96	5560	5.020	ng/uL	0.00
4) Pyridine-d5	4.023	84	20412	6.160	ng/u1	0.00
7) Phenol-d5	7.372	99	23051	6.044	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	71610	29.067	ng/u1	0.00
11) 2-Chlorophenol-d4	7.754	132	57803	21.870	ng/u1	-0.01
15) 4-Methylphenol-d8	8.929	113	40111	13.360	ng/u1	0.00
21) Nitrobenzene-d5	9.405	128	39797	28.711	ng/u1	0.00
24) 2-Nitrophenol-d4	10.127	143	41963	27.226	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.668	165	64130	24.701	ng/u1	0.00
31) 4-Chloroaniline-d4	11.185	131	91329	23.229	ng/u1	0.00
46) Dimethylphthalate-d6	14.246	166	277799	33.544	ng/u1	0.00
49) Acenaphthylene-d8	14.546	160	320869	31.098	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.045	143	7015	4.672	ng/u1	0.00
60) Fluorene-d10	15.838	176	238538	32.515	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.956	200	38325	26.118	ng/u1	0.00
73) Anthracene-d10	17.695	188	397921	34.775	ng/u1	-0.01
81) Pyrene-d10	19.969	212	460057	34.300	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.051	264	384812	34.353	ng/u1	-0.02

Target Compounds Qvalue

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

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