

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
 Data File : BG050847.D
 Acq On : 3 Nov 2021 12:49
 Operator : CG/JU
 Sample : M4412-05
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG361

Quant Time: Nov 03 14:29:31 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	25318	20.000	ng/u1	0.00
20) Naphthalene-d8	11.057	136	116296	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.852	164	83331	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.596	188	192804	20.000	ng/u1	0.00
79) Chrysene-d12	21.891	240	169596	20.000	ng/u1	-0.02
88) Perylene-d12	25.293	264	166833	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.589	96	3824	4.875	ng/uL	0.00
4) Pyridine-d5	4.018	84	20621	8.787	ng/u1	0.00
7) Phenol-d5	7.373	99	19467	7.207	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.543	67	58370	33.455	ng/u1	0.00
11) 2-Chlorophenol-d4	7.755	132	49138	26.251	ng/u1	0.00
15) 4-Methylphenol-d8	8.930	113	35387	16.642	ng/u1	0.00
21) Nitrobenzene-d5	9.400	128	33876	34.277	ng/u1	0.00
24) 2-Nitrophenol-d4	10.128	143	35943	32.708	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.669	165	54942	29.681	ng/u1	0.00
31) 4-Chloroaniline-d4	11.186	131	80653	28.771	ng/u1	0.00
46) Dimethylphthalate-d6	14.247	166	246000	38.586	ng/u1	0.00
49) Acenaphthylene-d8	14.553	160	276149	34.767	ng/u1	0.00
54) 4-Nitrophenol-d4	15.052	143	7327	6.338	ng/u1	0.00
60) Fluorene-d10	15.845	176	210731	37.313	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.957	200	36245	31.007	ng/u1	0.00
73) Anthracene-d10	17.696	188	360612	39.560	ng/u1	0.00
81) Pyrene-d10	19.970	212	424511	38.754	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.064	264	348433	37.780	ng/u1	0.00

Target Compounds	Qvalue

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

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