

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
 Data File : BG050839.D
 Acq On : 3 Nov 2021 7:21
 Operator : CG/JU
 Sample : M4400-04
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG333

Quant Time: Nov 03 08:12:11 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	27257	20.000	ng/ul	0.00
20) Naphthalene-d8	11.056	136	132039	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.852	164	94838	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.596	188	226925	20.000	ng/ul	-0.01
79) Chrysene-d12	21.897	240	205508	20.000	ng/ul	0.00
88) Perylene-d12	25.298	264	195058	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.589	96	3911	4.631	ng/uL	0.00
4) Pyridine-d5	4.017	84	18622	7.371	ng/ul	0.00
7) Phenol-d5	7.372	99	19736	6.787	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.543	67	57409	30.563	ng/ul	0.00
11) 2-Chlorophenol-d4	7.754	132	46984	23.315	ng/ul	-0.01
15) 4-Methylphenol-d8	8.929	113	36506	15.947	ng/ul	0.00
21) Nitrobenzene-d5	9.399	128	32430	28.902	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.128	143	34310	27.500	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.669	165	54269	25.822	ng/ul	0.00
31) 4-Chloroaniline-d4	11.186	131	78386	24.629	ng/ul	0.00
46) Dimethylphthalate-d6	14.247	166	253582	34.949	ng/ul	0.00
49) Acenaphthylene-d8	14.552	160	287300	31.782	ng/ul	0.00
54) 4-Nitrophenol-d4	15.046	143	8679	6.597	ng/ul	0.00
60) Fluorene-d10	15.845	176	216532	33.688	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.956	200	39529	28.732	ng/ul	0.00
73) Anthracene-d10	17.696	188	363066	33.840	ng/ul	-0.01
81) Pyrene-d10	19.975	212	441061	33.229	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.063	264	358969	33.290	ng/ul	0.00

Target Compounds Qvalue

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

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