

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122718\
 Data File : BG038851.D
 Acq On : 27 Dec 2018 19:58
 Operator : JU/SJ
 Sample : J6533-19
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COMP-5

Quant Time: Dec 28 01:33:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 15:26:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	23924	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	105193	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	76298	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	178049	20.00	ng	0.00
76) Chrysene-d12	21.72	240	186771	20.00	ng	0.00
87) Perylene-d12	25.02	264	202110	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	106764	79.15	ng	0.00
7) Phenol-d6	7.22	99	101301	54.71	ng	0.01
23) Nitrobenzene-d5	9.24	82	194538	94.48	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	190343	181.02	ng	0.00
45) 2-Fluorobiphenyl	13.31	172	543861	97.79	ng	0.00
79) Terphenyl-d14	20.04	244	871755	101.58	ng	0.00
Target Compounds						Qvalue
9) Benzaldehyde	7.20	77	2977	2.478	ng	96
10) Phenol	7.25	94	18434	9.370	ng	95
15) Benzyl Alcohol	8.31	79	186688	129.166	ng	100
17) 2-Methylphenol	8.50	107	3498	2.654	ng	# 90
20) 3+4-Methylphenols	8.83	107	17525	9.628	ng	95
27) 2,4-Dimethylphenol	10.04	122	7771	5.086	ng	95
32) Benzoic acid	10.31	122	136130	120.727	ng	# 82
50) Dimethylphthalate	14.14	163	17226	2.765	ng	99

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

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