Quantitation Report (Not Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG041619\
Data File : BG040523.D
Acq On : 17 Apr 2019 4:56
Operator : JU/SJ
Sample : K2411-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-14-D

Quant Time: Apr 17 07:00:10 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 28 05:47:00 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	2895	20.00	ng	-0.03
21) Naphthalene-d8	0.00	136	0	0.00	ng	-10.88
39) Acenaphthene-d10	0.00	164	0	0.00	ng	-14.70
64) Phenanthrene-d10	0.00	188	0	0.00	ng	-17.45
76) Chrysene-d12	0.00	240	0	0.00	ng	-21.73
87) Perylene-d12	0.00	264	0	0.00	ng	-25.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	387754	2226.62	ng	-0.02
7) Phenol-d6	7.20	99	466578	1938.66	ng	-0.03
23) Nitrobenzene-d5	9.22	82	389147	0.00	ng	-0.03
42) 2,4,6-Tribromophenol	16.16	330	168401	0.00	ng	-0.03
45) 2-Fluorobiphenyl	13.29	172	859776	0.00	ng	-0.03
79) Terphenyl-d14	20.02	244	1432207	0.00	ng	-0.03
Target Compounds						Qvalue
4) n-Nitrosodimethylamine	3.92	42	2305	22.602	ng	# 26
6) Aniline	7.58	93	726	2.322	ng	# 1
9) Benzaldehyde	7.20	77	766	4.640	ng	# 2
11) bis(2-Chloroethyl)ether	7.58	93	726	3.470	ng	# 1
15) Benzyl Alcohol	8.36	79	6034	31.132	ng	# 50
19) n-Nitroso-di-n-propylamine	9.22	70	50452	292.382	ng	# 73

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

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