

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092118\
 Data File : BG036875.D
 Acq On : 22 Sep 2018 1:53
 Operator : JU/SJ
 Sample : J4778-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SA-01-091918

Quant Time: Sep 22 04:30:54 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 20 20:07:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	1919	20.00	ng	0.00
21) Naphthalene-d8	0.00	136	0	0.00	ng	-10.86
38) Acenaphthene-d10	0.00	164	0	0.00	ng	-14.67
63) Phenanthrene-d10	0.00	188	0	0.00	ng	-17.42
75) Chrysene-d12	0.00	240	0	0.00	ng	-21.69
86) Perylene-d12	0.00	264	0	0.00	ng	-24.90

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.62	112	259083	2589.19	ng	0.00
7) Phenol-d6	7.21	99	325166	2100.35	ng	0.00
23) Nitrobenzene-d5	9.21	82	261638	0.00	ng	0.00
41) 2,4,6-Tribromophenol	16.16	330	170240	0.00	ng	0.00
44) 2-Fluorobiphenyl	13.29	172	585674	0.00	ng	0.00
78) Terphenyl-d14	20.03	244	1208716	0.00	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.54	88	123	2.686	ng	# 1
4) n-Nitrosodimethylamine	3.83	42	649	11.045	ng	# 1
6) Aniline	7.57	93	406	2.021	ng	# 1
9) Benzaldehyde	7.22	77	426	4.164	ng	# 4
10) Phenol	7.58	94	496	3.104	ng	# 1
11) bis(2-Chloroethyl)ether	7.57	93	406	2.747	ng	# 1
15) Benzyl Alcohol	8.37	79	4045	32.201	ng	# 59
16) 2,2'-oxybis(1-Chloropropan	8.57	45	674	2.375	ng	# 74
19) n-Nitroso-di-n-propylamine	9.21	70	35347	274.457	ng	# 71

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

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