

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG041619\
 Data File : BG040520.D
 Acq On : 17 Apr 2019 2:54
 Operator : JU/SJ
 Sample : PB118938BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB118938BL

Quant Time: Apr 17 04:58:59 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	2136	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	21.67	240	68	20.00	ng	-0.06
87) Perylene-d12	0.00	264	0	0.00	ng	-25.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	364975	2840.54	ng	-0.02
7) Phenol-d6	7.20	99	498250	2805.90	ng	-0.02
23) Nitrobenzene-d5	9.22	82	303315	0.00	ng	-0.03
42) 2,4,6-Tribromophenol	16.16	330	161250	0.00	ng	-0.03
45) 2-Fluorobiphenyl	13.30	172	729115	0.00	ng	-0.03
79) Terphenyl-d14	20.02	244	1232458	407735.72	ng	-0.03
Target Compounds						
2) 1,4-Dioxane	3.52	88	148	2.450	ng	# 1
4) n-Nitrosodimethylamine	3.84	42	234	3.110	ng	# 1
6) Aniline	7.57	93	507	2.198	ng	# 1
9) Benzaldehyde	7.20	77	1011	8.300	ng	# 2
11) bis(2-Chloroethyl)ether	7.57	93	507	3.284	ng	# 1
15) Benzyl Alcohol	8.36	79	4810	33.635	ng	# 57
19) n-Nitroso-di-n-propylamine	9.22	70	36809	289.117	ng	# 73
77) Benzidine	20.02	184	18918	7704.164	ng	# 94
78) Pyrene	20.02	202	4844	1083.241	ng	# 62
80) Butylbenzylphthalate	20.66	149	56	29.265	ng	# 23
81) Benzo(a)anthracene	22.11	228	198	47.273	ng	# 1
82) 3,3'-Dichlorobenzidine	21.92	252	75	47.072	ng	# 26
83) Chrysene	22.11	228	198	49.188	ng	# 7
84) Bis(2-ethylhexyl)phthalate	21.53	149	120	43.870	ng	# 50
85) Di-n-octyl phthalate	22.75	149	156	33.474	ng	# 1

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

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