

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041619\
 Data File : BF113802.D
 Acq On : 16 Apr 2019 19:22
 Operator : JU/SJ
 Sample : K2422-11
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP3-S

Quant Time: Apr 17 01:07:28 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 04 04:05:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	5273	20.00	ng	0.00
21) Naphthalene-d8	0.00	136	0	0.00	ng	-7.97
39) Acenaphthene-d10	0.00	164	0	0.00	ng	-9.72
64) Phenanthrene-d10	0.00	188	0	0.00	ng	-11.20
76) Chrysene-d12	13.84	240	244	20.00	ng	0.00
87) Perylene-d12	0.00	264	0	0.00	ng	-15.25
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	757800	2450.64	ng	0.00
7) Phenol-d6	6.33	99	1019939	2676.15	ng	0.00
23) Nitrobenzene-d5	7.24	82	587490	0.00	ng	-0.02
42) 2,4,6-Tribromophenol	10.50	330	221432	0.00	ng	-0.02
45) 2-Fluorobiphenyl	9.03	172	1089943	0.00	ng	-0.02
79) Terphenyl-d14	12.77	244	891131	74358.87	ng	-0.02
Target Compounds						
2) 1,4-Dioxane	2.40	88	771	5.264	ng	# 33
6) Aniline	6.45	93	986	2.134	ng	# 1
9) Benzaldehyde	6.20	77	1070	4.722	ng	# 1
10) Phenol	6.35	94	8110	18.632	ng	# 1
11) bis(2-Chloroethyl)ether	6.45	93	986	2.912	ng	# 1
15) Benzyl Alcohol	6.83	79	8523	33.074	ng	# 47
18) Hexachloroethane	7.21	117	642	4.530	ng	# 1
19) n-Nitroso-di-n-propylamine	7.24	70	72837	309.923	ng	# 77
77) Benzidine	12.76	184	9445	1040.936	ng	# 96
78) Pyrene	12.64	202	6254	336.858	ng	# 96
80) Butylbenzylphthalate	13.25	149	108	14.291	ng	# 25
81) Benzo(a)anthracene	13.83	228	3585	254.702	ng	# 87
83) Chrysene	13.87	228	3497	245.088	ng	# 90
84) Bis(2-ethylhexyl)phthalate	13.80	149	4130	451.996	ng	# 96
85) Di-n-octyl phthalate	14.44	149	867	58.374	ng	# 15
86) Indeno(1,2,3-cd)pyrene	16.64	276	2412	183.092	ng	# 87

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

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