

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041019\
 Data File : BF113708.D
 Acq On : 10 Apr 2019 20:52
 Operator : JU/SJ
 Sample : K2200-05 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-040919-C

Manual Integrations
 APPROVED

Sohil
 4/11/2019 2:39:52 PM

Quant Time: Apr 11 04:03:55 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	6.67	152	130829	20.00	ng	-0.01
21) Naphthalene-d8	7.96	136	486754	20.00	ng	-0.01
39) Acenaphthene-d10	9.71	164	205643	20.00	ng	-0.01
64) Phenanthrene-d10	11.19	188	343083	20.00	ng	-0.01
76) Chrysene-d12	13.83	240	327149	20.00	ng	-0.02
87) Perylene-d12	15.24	264	233066	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	319217	41.61	ng	0.00
7) Phenol-d6	6.33	99	430706	45.55	ng	0.00
23) Nitrobenzene-d5	7.25	82	240319	28.92	ng	-0.01
42) 2,4,6-Tribromophenol	10.50	330	89135	36.08	ng	-0.02
45) 2-Fluorobiphenyl	9.03	172	449827	35.22	ng	-0.02
79) Terphenyl-d14	12.77	244	394929	24.58	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	7.98	128	142256	6.032	ng	97
37) 2-Methylnaphthalene	8.67	142	50100	3.230	ng	95
38) 1-Methylnaphthalene	8.77	142	35865	2.398	ng	96
50) Dimethylphthalate	9.43	163	42614	2.811	ng	99
52) Acenaphthene	9.74	154	48790	4.080	ng	99
55) Dibenzofuran	9.92	168	71327	4.066	ng	96
58) Fluorene	10.26	166	98054	7.067	ng	99
71) Phenanthrene	11.22	178	542354	31.813	ng	99
72) Anthracene	11.27	178	193488	11.155	ng	97
73) Carbazole	11.43	167	45342	2.795	ng	98
75) Fluoranthene	12.40	202	733266	40.139	ng	99
78) Pyrene	12.63	202	666689	26.783	ng	99
81) Benzo(a)anthracene	13.82	228	345195	18.292	ng	98
83) Chrysene	13.85	228	282930	14.789	ng	97
86) Indeno(1,2,3-cd)pyrene	16.60	276	104970	5.943	ng	96
88) Benzo(b)fluoranthene	14.84	252	280093m	19.633	ng	
89) Benzo(k)fluoranthene	14.86	252	95562m	7.464	ng	
90) Benzo(a)pyrene	15.18	252	200227	15.794	ng	98
91) Dibenzo(a,h)anthracene	16.60	278	27607	2.329	ng	# 75
92) Benzo(a,h,i)perylene	17.02	276	100592	8.313	ng	96

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

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