

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041019\
 Data File : BF113707.D
 Acq On : 10 Apr 2019 20:23
 Operator : JU/SJ
 Sample : K2200-03 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 4/11/2019 2:39:49 PM

Quant Time: Apr 11 04:00:40 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	136559	20.00	ng	-0.01
21) Naphthalene-d8	7.96	136	522775	20.00	ng	-0.01
39) Acenaphthene-d10	9.71	164	227008	20.00	ng	-0.01
64) Phenanthrene-d10	11.19	188	370319	20.00	ng	-0.01
76) Chrysene-d12	13.83	240	338942	20.00	ng	-0.02
87) Perylene-d12	15.24	264	240778	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	407804	50.92	ng	0.00
7) Phenol-d6	6.33	99	559022	56.64	ng	0.00
23) Nitrobenzene-d5	7.25	82	307416	34.44	ng	-0.01
42) 2,4,6-Tribromophenol	10.50	330	122389	44.88	ng	-0.02
45) 2-Fluorobiphenyl	9.03	172	590472	41.87	ng	-0.02
79) Terphenyl-d14	12.77	244	504931	30.33	ng	-0.02
Target Compounds						Qvalue
49) Acenaphthylene	9.57	152	51906	2.251	ng	98
50) Dimethylphthalate	9.43	163	51964	3.105	ng	99
58) Fluorene	10.26	166	55460	3.621	ng	97
71) Phenanthrene	11.22	178	421755	22.920	ng	98
72) Anthracene	11.27	178	101432	5.418	ng	98
75) Fluoranthene	12.40	202	404578	20.518	ng	99
78) Pyrene	12.63	202	493336	19.129	ng	99
81) Benzo(a)anthracene	13.82	228	220988	11.303	ng	95
83) Chrysene	13.85	228	208591	10.524	ng	98
86) Indeno(1,2,3-cd)pyrene	16.60	276	52300	2.858	ng	99
88) Benzo(b)fluoranthene	14.84	252	148660m	10.086	ng	
89) Benzo(k)fluoranthene	14.86	252	49411m	3.735	ng	
90) Benzo(a)pyrene	15.18	252	120373	9.191	ng	# 95
92) Benzo(a,h,i)perylene	17.02	276	59320	4.745	ng	94

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

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