

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041019\
 Data File : BF113710.D
 Acq On : 10 Apr 2019 21:50
 Operator : JU/SJ
 Sample : K2203-07 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-040919-B

Manual Integrations
 APPROVED

Sohil
 4/11/2019 2:39:58 PM

Quant Time: Apr 11 04:19:59 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	127856	20.00	ng	-0.01
21) Naphthalene-d8	7.96	136	460982	20.00	ng	0.00
39) Acenaphthene-d10	9.71	164	184341	20.00	ng	-0.01
64) Phenanthrene-d10	11.19	188	353592	20.00	ng	-0.01
76) Chrysene-d12	13.83	240	313342	20.00	ng	-0.02
87) Perylene-d12	15.24	264	219587	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	341419	45.54	ng	0.00
7) Phenol-d6	6.33	99	463086	50.11	ng	0.00
23) Nitrobenzene-d5	7.25	82	249035	31.64	ng	-0.01
42) 2,4,6-Tribromophenol	10.50	330	97828	44.18	ng	-0.02
45) 2-Fluorobiphenyl	9.03	172	462166	40.36	ng	-0.02
79) Terphenyl-d14	12.77	244	469551	30.51	ng	-0.02
Target Compounds						Qvalue
49) Acenaphthylene	9.57	152	63158	3.372	ng	98
50) Dimethylphthalate	9.43	163	40316	2.967	ng	100
71) Phenanthrene	11.22	178	182810	10.405	ng	97
72) Anthracene	11.27	178	93349	5.222	ng	99
75) Fluoranthene	12.40	202	877929	46.629	ng	100
78) Pyrene	12.63	202	854955	35.859	ng	99
81) Benzo(a)anthracene	13.82	228	467541	25.866	ng	97
83) Chrysene	13.85	228	419537	22.896	ng	99
86) Indeno(1,2,3-cd)pyrene	16.60	276	181881	10.751	ng	95
88) Benzo(b)fluoranthene	14.84	252	432346m	32.165	ng	
89) Benzo(k)fluoranthene	14.86	252	139285m	11.546	ng	
90) Benzo(a)pyrene	15.18	252	295072	24.704	ng	99
91) Dibenzo(a,h)anthracene	16.60	278	45560	4.079	ng	# 78
92) Benzo(g,h,i)perylene	17.01	276	179231	15.721	ng	97

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

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