

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041619\
 Data File : BF113806.D
 Acq On : 16 Apr 2019 21:18
 Operator : JU/SJ
 Sample : K2203-09 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-01-041519-A

Manual Integrations
 APPROVED

Sohil
 4/17/2019 3:45:52 PM

Quant Time: Apr 17 04:38:06 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	59767	20.00	ng	-0.01
21) Naphthalene-d8	7.96	136	214886	20.00	ng	-0.01
39) Acenaphthene-d10	9.70	164	88153	20.00	ng	-0.02
64) Phenanthrene-d10	11.19	188	162422	20.00	ng	-0.02
76) Chrysene-d12	13.83	240	168321	20.00	ng	-0.02
87) Perylene-d12	15.24	264	155388	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	339999	97.01	ng	0.00
7) Phenol-d6	6.33	99	455777	105.51	ng	0.00
23) Nitrobenzene-d5	7.24	82	248153	67.64	ng	-0.02
42) 2,4,6-Tribromophenol	10.50	330	93460	88.26	ng	-0.01
45) 2-Fluorobiphenyl	9.03	172	467665	85.41	ng	-0.02
79) Terphenyl-d14	12.76	244	454370	54.96	ng	-0.03
Target Compounds						
50) Dimethylphthalate	9.43	163	51658	7.950	ng	Qvalue 100
52) Acenaphthene	9.74	154	12501	2.438	ng	93
58) Fluorene	10.26	166	22439	3.773	ng	96
71) Phenanthrene	11.22	178	178484	22.115	ng	97
72) Anthracene	11.27	178	71077	8.656	ng	98
75) Fluoranthene	12.40	202	390270	45.126	ng	99
78) Pyrene	12.63	202	358716	28.009	ng	99
81) Benzo(a)anthracene	13.82	228	196268	20.214	ng	99
83) Chrysene	13.85	228	169801	17.251	ng	99
86) Indeno(1,2,3-cd)pyrene	16.62	276	61855	6.806	ng	93
88) Benzo(b)fluoranthene	14.84	252	198561m	20.876	ng	
89) Benzo(k)fluoranthene	14.87	252	52828m	6.188	ng	
90) Benzo(a)pyrene	15.19	252	133449	15.789	ng	99
92) Benzo(g,h,i)perylene	17.03	276	53853	6.675	ng	94

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

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