

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042519\
 Data File : BF113998.D
 Acq On : 26 Apr 2019 00:24
 Operator : JU/SJ
 Sample : K2203-17 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-03-042419-A

Manual Integrations
 APPROVED

Sohil
 4/29/2019 5:49:59 AM

Quant Time: Apr 26 06:49:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	112568	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	459887	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	251634	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	525191	20.00	ng	0.00
76) Chrysene-d12	14.05	240	380470	20.00	ng	-0.02
87) Perylene-d12	15.54	264	407591	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	123117	18.71	ng	0.00
7) Phenol-d6	6.51	99	169866	20.57	ng	-0.01
23) Nitrobenzene-d5	7.45	82	96474	12.95	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	59749	19.49	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	211156	13.12	ng	0.00
79) Terphenyl-d14	12.99	244	263542	14.20	ng	-0.01
Target Compounds						
71) Phenanthrene	11.44	178	115491	4.376	ng	98
75) Fluoranthene	12.63	202	215689	7.490	ng	97
78) Pyrene	12.86	202	179685	6.753	ng	97
81) Benzo(a)anthracene	14.04	228	87999	3.926	ng	99
83) Chrysene	14.08	228	74364	3.406	ng	98
86) Indeno(1,2,3-cd)pyrene	17.02	276	42953	2.077	ng	95
88) Benzo(b)fluoranthene	15.10	252	90876m	3.524	ng	
90) Benzo(a)pyrene	15.47	252	68926	3.090	ng	99

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

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