

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040419\
 Data File : BF113598.D
 Acq On : 4 Apr 2019 22:32
 Operator : JU/SJ
 Sample : K2221-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S2B(2-10)COMP

Manual Integrations
 APPROVED

Sohil
 4/6/2019 12:38:52 AM

Quant Time: Apr 05 05:25:59 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	6.68	152	145563	20.00	ng	0.00
21) Naphthalene-d8	7.96	136	553097	20.00	ng	0.00
39) Acenaphthene-d10	9.71	164	280168	20.00	ng	-0.01
64) Phenanthrene-d10	11.19	188	569315	20.00	ng	-0.02
76) Chrysene-d12	13.82	240	425353	20.00	ng	-0.03
87) Perylene-d12	15.22	264	384311	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	805161	94.32	ng	0.00
7) Phenol-d6	6.33	99	1066915	101.41	ng	-0.01
23) Nitrobenzene-d5	7.25	82	647369	68.56	ng	-0.01
42) 2,4,6-Tribromophenol	10.50	330	332117	98.68	ng	-0.02
45) 2-Fluorobiphenyl	9.03	172	1259739	72.39	ng	-0.02
79) Terphenyl-d14	12.77	244	1500666	71.83	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.43	163	154677	7.490	ng	Qvalue 100
71) Phenanthrene	11.21	178	115473	4.082	ng	97
75) Fluoranthene	12.40	202	703111	23.194	ng	98
78) Pyrene	12.62	202	605812	18.718	ng	99
81) Benzo(a)anthracene	13.80	228	260422	10.614	ng	99
83) Chrysene	13.84	228	330559	13.290	ng	99
84) Bis(2-ethylhexyl)phthalate	13.80	149	48317	3.033	ng	99
86) Indeno(1,2,3-cd)pyrene	16.56	276	81398	3.544	ng	98
88) Benzo(b)fluoranthene	14.82	252	259066m	11.013	ng	
89) Benzo(k)fluoranthene	14.84	252	81222m	3.847	ng	
90) Benzo(a)pyrene	15.16	252	110034	5.264	ng	98
92) Benzo(g,h,i)perylene	16.97	276	77603	3.889	ng	98

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

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