

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
 Data File : BF113519.D
 Acq On : 2 Apr 2019 21:41
 Operator : JU/SJ
 Sample : K2182-09
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S7C(12-20)COMP

Manual Integrations
 APPROVED

Sohil
 4/3/2019 12:42:44 PM

Quant Time: Apr 03 07:42:45 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	150511	20.00	ng	-0.01
21) Naphthalene-d8	7.97	136	554817	20.00	ng	-0.01
39) Acenaphthene-d10	9.72	164	261710	20.00	ng	-0.02
64) Phenanthrene-d10	11.20	188	461479	20.00	ng	-0.02
76) Chrysene-d12	13.83	240	357419	20.00	ng	-0.02
87) Perylene-d12	15.23	264	399390	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	852272	92.59	ng	0.00
7) Phenol-d6	6.35	99	1102196	94.65	ng	-0.01
23) Nitrobenzene-d5	7.26	82	654771	70.05	ng	-0.01
42) 2,4,6-Tribromophenol	10.51	330	297088	91.90	ng	-0.02
45) 2-Fluorobiphenyl	9.05	172	1233209	70.94	ng	-0.02
79) Terphenyl-d14	12.78	244	1068780	65.91	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.44	163	87293	4.501	ng	Qvalue 100
71) Phenanthrene	11.22	178	64773	2.826	ng	97
75) Fluoranthene	12.40	202	162170	6.260	ng	99
78) Pyrene	12.63	202	160721	6.489	ng	98
81) Benzo(a)anthracene	13.82	228	108270	5.292	ng	99
83) Chrysene	13.85	228	110117	5.299	ng	99
86) Indeno(1,2,3-cd)pyrene	16.58	276	45700	2.562	ng	96
88) Benzo(b)fluoranthene	14.84	252	173793m	7.005	ng	
90) Benzo(a)pyrene	15.17	252	66612	3.029	ng	99
92) Benzo(g,h,i)perylene	16.99	276	38438	2.005	ng	# 94

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

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