

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031419\
 Data File : BF113053.D
 Acq On : 14 Mar 2019 16:50
 Operator : JU/SJ
 Sample : K1866-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MR-MH-03-COMP

Manual Integrations
 APPROVED

Sohil
 3/15/2019 10:03:27 AM

Quant Time: Mar 15 07:21:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 13 03:42:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	214021	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	815217	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	382444	20.00	ng	-0.01
64) Phenanthrene-d10	11.23	188	803832	20.00	ng	-0.01
76) Chrysene-d12	13.86	240	730902	20.00	ng	0.00
87) Perylene-d12	15.27	264	706788	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1207853	87.79	ng	0.00
7) Phenol-d6	6.37	99	1566289	90.63	ng	0.00
23) Nitrobenzene-d5	7.29	82	972802	71.40	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	490884	120.90	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1712354	71.04	ng	0.00
79) Terphenyl-d14	12.82	244	1948744	51.68	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.47	163	186497	6.613	ng	98
71) Phenanthrene	11.26	178	132152	3.304	ng	98
75) Fluoranthene	12.44	202	432595	10.096	ng	96
78) Pyrene	12.67	202	431402	7.001	ng	99
81) Benzo(a)anthracene	13.85	228	230465	4.550	ng	97
83) Chrysene	13.89	228	219780	4.429	ng	98
86) Indeno(1,2,3-cd)pyrene	16.62	276	149186	3.527	ng	# 89
88) Benzo(b)fluoranthene	14.87	252	293753m	6.425	ng	
89) Benzo(k)fluoranthene	14.89	252	86588m	2.091	ng	
90) Benzo(a)pyrene	15.21	252	207687	5.136	ng	97
92) Benzo(g,h,i)perylene	17.03	276	147925	4.269	ng	96

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

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