

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031519\
 Data File : BF113061.D
 Acq On : 15 Mar 2019 13:07
 Operator : JU/SJ
 Sample : K1905-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RT5124

Manual Integrations
 APPROVED

Sohil
 3/18/2019 2:59:26 PM

Quant Time: Mar 15 23:04:49 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	164245	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	641772	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	311688	20.00	ng	-0.01
64) Phenanthrene-d10	11.24	188	640459	20.00	ng	0.00
76) Chrysene-d12	13.86	240	609631	20.00	ng	0.00
87) Perylene-d12	15.27	264	583258	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1122016	106.26	ng	0.00
7) Phenol-d6	6.37	99	1527081	115.13	ng	0.00
23) Nitrobenzene-d5	7.29	82	888275	82.82	ng	-0.01
42) 2,4,6-Tribromophenol	10.54	330	475777	143.78	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1524524	79.72	ng	0.00
79) Terphenyl-d14	12.82	244	1741626	55.38	ng	0.00
Target Compounds						
10) Phenol	6.39	94	31640	2.044	ng	96
50) Dimethylphthalate	9.47	163	129647	5.641	ng	# 98
71) Phenanthrene	11.26	178	112875	3.542	ng	# 94
75) Fluoranthene	12.44	202	380327	11.140	ng	95
78) Pyrene	12.67	202	314727	6.124	ng	100
81) Benzo(a)anthracene	13.85	228	127936	3.028	ng	96
83) Chrysene	13.89	228	171542	4.145	ng	98
86) Indeno(1,2,3-cd)pyrene	16.62	276	106978	3.032	ng	# 90
88) Benzo(b)fluoranthene	14.87	252	253638m	6.723	ng	
89) Benzo(k)fluoranthene	14.89	252	69425m	2.032	ng	
90) Benzo(a)pyrene	15.21	252	134169	4.021	ng	98
92) Benzo(a,h,i)perylene	17.03	276	105093	3.675	ng	96

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

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