

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031219\
 Data File : BF113020.D
 Acq On : 12 Mar 2019 23:23
 Operator : JU/SJ
 Sample : K1847-14 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-7B-C-1-031119

Manual Integrations
 APPROVED

Sohil
 3/13/2019 11:40:51 AM

Quant Time: Mar 13 04:48:16 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	190878	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	696032	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	288750	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	511451	20.00	ng	0.00
76) Chrysene-d12	13.86	240	530948	20.00	ng	0.00
87) Perylene-d12	15.27	264	405479	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	680577	55.46	ng	0.00
7) Phenol-d6	6.37	99	877990	56.96	ng	0.00
23) Nitrobenzene-d5	7.29	82	510999	43.93	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	204928	66.85	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	915303	46.55	ng	0.00
79) Terphenyl-d14	12.82	244	838049	30.60	ng	0.00
Target Compounds						
10) Phenol	6.38	94	64063	3.562	ng	86
50) Dimethylphthalate	9.48	163	87087	4.090	ng	99
71) Phenanthrene	11.26	178	149309	5.868	ng	98
75) Fluoranthene	12.44	202	359465	13.185	ng	95
78) Pyrene	12.67	202	312018	6.971	ng	97
81) Benzo(a)anthracene	13.86	228	180678	4.910	ng	99
83) Chrysene	13.89	228	181544	5.037	ng	98
88) Benzo(b)fluoranthene	14.87	252	181475m	6.919	ng	
89) Benzo(k)fluoranthene	14.89	252	79869m	3.362	ng	
90) Benzo(a)pyrene	15.21	252	120757	5.205	ng	98
92) Benzo(g,h,i)perylene	17.03	276	60622	3.050	ng	97

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

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