

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031219\
 Data File : BF113019.D
 Acq On : 12 Mar 2019 22:55
 Operator : JU/SJ
 Sample : K1847-18
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-7B-C-5-031119

Manual Integrations
 APPROVED

Sohil
 3/13/2019 11:40:49 AM

Quant Time: Mar 13 04:46:32 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	174823	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	643590	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	300818	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	546962	20.00	ng	0.00
76) Chrysene-d12	13.86	240	476326	20.00	ng	0.00
87) Perylene-d12	15.27	264	391296	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	930093	82.76	ng	0.00
7) Phenol-d6	6.37	99	1231119	87.20	ng	0.00
23) Nitrobenzene-d5	7.29	82	725510	67.45	ng	0.00
42) 2,4,6-Tribromophenol	10.55	330	345257	108.11	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1336579	70.34	ng	0.00
79) Terphenyl-d14	12.82	244	1235514	50.28	ng	0.00
Target Compounds						
10) Phenol	6.38	94	86553	5.254	ng	91
50) Dimethylphthalate	9.48	163	146351	6.598	ng	99
71) Phenanthrene	11.26	178	115967	4.261	ng	97
75) Fluoranthene	12.45	202	156820	5.379	ng	97
78) Pyrene	12.67	202	233066	5.804	ng	99
81) Benzo(a)anthracene	13.86	228	92182m	2.792	ng	
83) Chrysene	13.89	228	102291	3.163	ng	96
88) Benzo(b)fluoranthene	14.87	252	73108m	2.888	ng	
90) Benzo(a)pyrene	15.21	252	51453	2.298	ng	99

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

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