

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022719\
 Data File : BF112734.D
 Acq On : 27 Feb 2019 22:20
 Operator : JU/SJ
 Sample : K1350-11 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-022619-A

Manual Integrations
 APPROVED

Sohil
 2/28/2019 12:47:22 PM

Quant Time: Feb 28 03:27:04 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 27 16:08:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	175043	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	608471	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	244617	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	492461	20.00	ng	0.00
76) Chrysene-d12	13.87	240	491729	20.00	ng	0.00
87) Perylene-d12	15.28	264	335772	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	225429	20.03	ng	0.00
7) Phenol-d6	6.36	99	279502	19.77	ng	-0.01
23) Nitrobenzene-d5	7.29	82	140826	13.85	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	50744	19.54	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	233541	3.79	ng	0.00
79) Terphenyl-d14	12.83	244	263306	10.38	ng	0.00
Target Compounds						
75) Fluoranthene	12.45	202	167622	6.385	ng	98
78) Pyrene	12.67	202	192947	4.654	ng	98
81) Benzo(a)anthracene	13.86	228	124001	3.639	ng	96
83) Chrysene	13.89	228	122031	3.656	ng	96
88) Benzo(b)fluoranthene	14.88	252	140458m	6.467	ng	
89) Benzo(k)fluoranthene	14.90	252	51509m	2.618	ng	
90) Benzo(a)pyrene	15.22	252	97702	5.086	ng	94
92) Benzo(g,h,i)perylene	17.03	276	52548	3.192	ng	100

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

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