

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032819\
 Data File : BF113373.D
 Acq On : 28 Mar 2019 14:42
 Operator : JU/SJ
 Sample : K1950-03 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-SF-03-005-B

Manual Integrations
 APPROVED

Sohil
 3/29/2019 4:47:18 PM

Quant Time: Mar 28 15:09:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	166880	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	596985	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	251841	20.00	ng	0.00
64) Phenanthrene-d10	11.21	188	482653	20.00	ng	0.00
76) Chrysene-d12	13.84	240	394660	20.00	ng	0.00
87) Perylene-d12	15.25	264	296306	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	202041	19.80	ng	0.00
7) Phenol-d6	6.35	99	345156	26.73	ng	-0.01
23) Nitrobenzene-d5	7.26	82	183116	18.21	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	31043	9.98	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	389029	8.30	ng	0.00
79) Terphenyl-d14	12.79	244	399191	22.29	ng	0.00
Target Compounds						
71) Phenanthrene	11.23	178	191355	7.984	ng	99
75) Fluoranthene	12.42	202	215105	7.939	ng	98
78) Pyrene	12.64	202	209009	7.642	ng	98
81) Benzo(a)anthracene	13.83	228	81747	3.619	ng	# 90
83) Chrysene	13.86	228	86440	3.767	ng	# 93
88) Benzo(b)fluoranthene	14.85	252	70239m	3.816	ng	
90) Benzo(a)pyrene	15.19	252	49473	3.032	ng	# 71
92) Benzo(g,h,i)perylene	17.02	276	29746	2.091	ng	# 72

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

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