

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031819\
 Data File : BF113093.D
 Acq On : 18 Mar 2019 16:32
 Operator : JU/SJ
 Sample : K1933-08 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-8

Manual Integrations
 APPROVED

Sohil
 3/19/2019 5:09:46 PM

Quant Time: Mar 19 05:28:06 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	176647	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	667586	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	339246	20.00	ng	-0.01
64) Phenanthrene-d10	11.24	188	601044	20.00	ng	0.00
76) Chrysene-d12	13.87	240	489208	20.00	ng	0.00
87) Perylene-d12	15.28	264	506288	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	414398	36.49	ng	0.00
7) Phenol-d6	6.37	99	603204	42.29	ng	0.00
23) Nitrobenzene-d5	7.29	82	336881	30.20	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	162348	45.08	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	737975	30.10	ng	-0.01
79) Terphenyl-d14	12.82	244	591291	23.43	ng	0.00
Target Compounds						
75) Fluoranthene	12.45	202	83055	2.592	ng	95
78) Pyrene	12.67	202	87173	2.114	ng	99
81) Benzo(a)anthracene	13.86	228	93839	2.768	ng	93
83) Chrysene	13.89	228	117374	3.534	ng	97
84) Bis(2-ethylhexyl)phthalate	13.86	149	75447	3.533	ng	100
88) Benzo(b)fluoranthene	14.87	252	125672m	3.838	ng	
90) Benzo(a)pyrene	15.22	252	128903	4.450	ng	95
92) Benzo(g,h,i)perylene	17.04	276	74115	2.986	ng	96

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

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