

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082318\
 Data File : BF108442.D
 Acq On : 24 Aug 2018 1:05
 Operator : JU/SJ
 Sample : J4283-03 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-082118

Manual Integrations
 APPROVED

Sohil
 8/24/2018 12:14:30 PM

Quant Time: Aug 24 10:57:41 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 22 11:01:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	131503	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	495720	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	223179	20.00	ng	-0.01
63) Phenanthrene-d10	11.54	188	356691	20.00	ng	0.00
75) Chrysene-d12	14.19	240	269322	20.00	ng	0.00
86) Perylene-d12	15.75	264	237428	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	100989	13.90	ng	-0.01
7) Phenol-d6	6.61	99	140299	14.54	ng	-0.02
23) Nitrobenzene-d5	7.56	82	87318	9.83	ng	-0.01
41) 2,4,6-Tribromophenol	10.84	330	22411	10.07	ng	-0.01
44) 2-Fluorobiphenyl	9.35	172	174426	12.55	ng	-0.01
78) Terphenyl-d14	13.12	244	123719	11.68	ng	-0.01
Target Compounds						
57) Fluorene	10.60	166	39905	2.763	ng	Qvalue 97
70) Phenanthrene	11.57	178	294034	17.477	ng	98
71) Anthracene	11.61	178	93395	5.416	ng	98
74) Fluoranthene	12.76	202	454573	24.757	ng	94
77) Pyrene	13.00	202	406318	23.830	ng	99
80) Benzo(a)anthracene	14.18	228	233424	15.102	ng	100
82) Chrysene	14.22	228	183473	12.948	ng	99
85) Indeno(1,2,3-cd)pyrene	17.35	276	94270m	7.678	ng	
87) Benzo(b)fluoranthene	15.28	252	278033m	19.597	ng	
88) Benzo(k)fluoranthene	15.31	252	83072m	6.370	ng	
89) Benzo(a)pyrene	15.68	252	198392	15.616	ng	98
90) Dibenzo(a,h)anthracene	17.35	278	23487m	2.234	ng	
91) Benzo(a,h,i)perylene	17.85	276	85731	8.283	ng	# 88

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

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