

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
 Data File : BF106361.D
 Acq On : 12 Jun 2018 22:44
 Operator : JU/SJ
 Sample : J3322-04 10X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB02-3_E_5

Manual Integrations
 APPROVED

Sohil
 6/14/2018 3:53:22 PM

Quant Time: Jun 13 15:06:08 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 11 15:00:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	127901	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	458303	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	197565	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	410427	20.00	ng	0.00
75) Chrysene-d12	14.17	240	317790	20.00	ng	0.01
86) Perylene-d12	15.71	264	279448	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	88863	11.76	ng	0.06
7) Phenol-d6	6.66	99	108240	11.34	ng	0.07
23) Nitrobenzene-d5	7.53	82	60320	7.74	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	20647	10.86	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	109992	0.84	ng	0.00
78) Terphenyl-d14	13.09	244	147325	11.51	ng	0.00
Target Compounds						
31) Naphthalene	8.28	128	1174170	56.675	ng	98
37) 2-Methylnaphthalene	8.97	142	135851	9.627	ng	98
45) 1,1'-Biphenyl	9.43	154	101251	6.617	ng	98
48) Acenaphthylene	9.88	152	1056851	56.383	ng	98
51) Acenaphthene	10.05	154	281997	23.296	ng	99
54) Dibenzofuran	10.22	168	258007	15.828	ng	97
57) Fluorene	10.57	166	418426	32.399	ng	100
70) Phenanthrene	11.59	178	894137	44.484	ng	99
71) Anthracene	11.59	178	894137	44.403	ng	98
72) Carbazole	11.74	167	181423	9.759	ng	99
74) Fluoranthene	12.74	202	2645413	123.411	ng	93
77) Pylene	12.98	202	2695642	135.489	ng	# 87
80) Benzo(a)anthracene	14.16	228	2069575	109.435	ng	# 74
82) Chrysene	14.20	228	1708142	98.930	ng	# 90
85) Indeno(1,2,3-cd)pyrene	17.30	276	1210568	87.850	ng	# 89
87) Benzo(b)fluoranthene	15.27	252	2640573m	156.318	ng	
88) Benzo(k)fluoranthene	15.29	252	646538m	38.646	ng	
89) Benzo(a)pyrene	15.66	252	1869390	123.035	ng	93
90) Dibenzo(a,h)anthracene	17.31	278	318414m	25.137	ng	
91) Benzo(a,h,i)perylene	17.79	276	1213405	98.568	ng	93

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

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