

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051119\
 Data File : BF114185.D
 Acq On : 12 May 2019 5:42
 Operator : JU/SJ
 Sample : K2765-08 10X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P026-SS004-0002-01

Manual Integrations
 APPROVED

Sohil
 5/13/2019 8:36:19 PM

Quant Time: May 13 08:16:51 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 10 15:56:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	252618	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	946827	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	415006	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	745325	20.00	ng	0.00
76) Chrysene-d12	14.01	240	614525	20.00	ng	-0.01
87) Perylene-d12	15.48	264	659085	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	201205	12.82	ng	0.00
7) Phenol-d6	6.48	99	273912	14.75	ng	-0.01
23) Nitrobenzene-d5	7.40	82	149676	8.87	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	51770	12.07	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	270714	9.87	ng	0.00
79) Terphenyl-d14	12.94	244	228697	7.67	ng	-0.01
Target Compounds						
49) Acenaphthylene	9.75	152	318226	7.754	ng	97
71) Phenanthrene	11.39	178	1530453	42.207	ng	99
72) Anthracene	11.45	178	198904	5.461	ng	97
75) Fluoranthene	12.59	202	1906962	48.836	ng	97
78) Pyrene	12.82	202	3228664	65.311	ng	97
81) Benzo(a)anthracene	14.00	228	1299261m	32.688	ng	
83) Chrysene	14.04	228	1376771	35.335	ng	99
86) Indeno(1,2,3-cd)pyrene	16.94	276	474628	13.783	ng	97
88) Benzo(b)fluoranthene	15.06	252	1249735m	29.597	ng	
89) Benzo(k)fluoranthene	15.07	252	400465m	10.325	ng	
90) Benzo(a)pyrene	15.42	252	930187	25.031	ng	98
91) Dibenzo(a,h)anthracene	16.94	278	127766m	4.138	ng	
92) Benzo(a,h,i)perylene	17.39	276	524358	16.944	ng	98

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

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