

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051119\
 Data File : BF114169.D
 Acq On : 11 May 2019 22:41
 Operator : JU/SJ
 Sample : K2765-09
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P026-SS004-0206-01

Manual Integrations
 APPROVED

Sohil
 5/13/2019 8:35:59 PM

Quant Time: May 13 08:11:50 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	245035	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	914409	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	434523	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	879071	20.00	ng	0.00
76) Chrysene-d12	14.00	240	566490	20.00	ng	-0.02
87) Perylene-d12	15.47	264	684611	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1671833	109.80	ng	0.00
7) Phenol-d6	6.49	99	2236360	124.18	ng	0.00
23) Nitrobenzene-d5	7.41	82	1351636	82.95	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	563250	125.38	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	2259887	78.67	ng	0.00
79) Terphenyl-d14	12.95	244	2251479	81.93	ng	0.00
Target Compounds						
49) Acenaphthylene	9.74	152	219038	5.098	ng	97
50) Dimethylphthalate	9.59	163	397539	12.463	ng	100
71) Phenanthrene	11.39	178	382294	8.939	ng	97
72) Anthracene	11.44	178	97422	2.268	ng	97
75) Fluoranthene	12.58	202	927111	20.130	ng	100
78) Pyrene	12.81	202	1615779	35.456	ng	99
81) Benzo(a)anthracene	13.99	228	615563m	16.800	ng	
83) Chrysene	14.03	228	781498	21.758	ng	99
86) Indeno(1,2,3-cd)pyrene	16.93	276	635654	20.025	ng	96
88) Benzo(b)fluoranthene	15.04	252	1073392m	24.473	ng	
89) Benzo(k)fluoranthene	15.07	252	237135m	5.886	ng	
90) Benzo(a)pyrene	15.40	252	661116	17.127	ng	99
91) Dibenzo(a,h)anthracene	16.93	278	163047m	5.083	ng	
92) Benzo(a,h,i)perylene	17.37	276	750274	23.341	ng	99

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

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