

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021220\
 Data File : BF118911.D
 Acq On : 12 Feb 2020 11:59
 Operator : CG/JU
 Sample : L1482-03DL 50X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP6-EP1-1.5DL

Manual Integrations
 APPROVED

mohammad
 2/13/2020 3:34:18 PM

Quant Time: Feb 13 00:21:28 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF020320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 03 16:26:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	205192	20.00	ng	-0.01
21) Naphthalene-d8	8.09	136	754496	20.00	ng	-0.02
39) Acenaphthene-d10	9.84	164	382953	20.00	ng	-0.02
64) Phenanthrene-d10	11.33	188	636470	20.00	ng	-0.01
76) Chrysene-d12	13.96	240	434189	20.00	ng	-0.02
87) Perylene-d12	15.42	264	513926	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	16252	1.50	ng	0.00
7) Phenol-d6	6.45	99	22073	1.64	ng	0.00
23) Nitrobenzene-d5	7.36	82	15479	1.22	ng	-0.02
42) 2,4,6-Tribromophenol	10.63	330	7791	1.70	ng	-0.02
45) 2-Fluorobiphenyl	9.16	172	38257	1.79	ng	-0.02
79) Terphenyl-d14	12.90	244	40686	1.92	ng	-0.02
Target Compounds						
31) Naphthalene	8.12	128	3117562	90.344	ng	99
37) 2-Methylnaphthalene	8.80	142	1338577	55.931	ng	96
38) 1-Methylnaphthalene	8.90	142	1024390	45.501	ng	100
46) 1,1'-Biphenyl	9.26	154	333902	12.605	ng	97
49) Acenaphthylene	9.70	152	155544	4.892	ng	# 90
52) Acenaphthene	9.87	154	130860	6.245	ng	93
55) Dibenzofuran	10.05	168	69088	2.282	ng	# 86
58) Fluorene	10.39	166	678873	29.800	ng	98
71) Phenanthrene	11.36	178	2467533	78.796	ng	100
72) Anthracene	11.40	178	353696	11.353	ng	97
73) Carbazole	11.56	167	65244	2.386	ng	99
75) Fluoranthene	12.54	202	936002	27.376	ng	99
78) Pylene	12.77	202	1417583	47.591	ng	99
81) Benzo(a)anthracene	13.95	228	510374	19.754	ng	97
83) Chrysene	13.99	228	515770	19.890	ng	97
86) Indeno(1,2,3-cd)pyrene	16.85	276	213578	8.198	ng	98
88) Benzo(b)fluoranthene	15.00	252	381627m	11.967	ng	
89) Benzo(k)fluoranthene	15.01	252	138508m	4.988	ng	
90) Benzo(a)pyrene	15.35	252	465838	17.069	ng	99
91) Dibenz(a,h)anthracene	16.85	278	73185	3.021	ng	# 79
92) Benzo(g,h,i)perylene	17.28	276	240832	10.336	ng	99

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

