

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021220\
 Data File : BF118912.D
 Acq On : 12 Feb 2020 12:27
 Operator : CG/JU
 Sample : L1482-02DL 25X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP7-EP2-2DL

Manual Integrations
 APPROVED

mohammad
 2/13/2020 3:34:21 PM

Quant Time: Feb 13 00:26:44 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	265961	20.00	ng	-0.01
21) Naphthalene-d8	8.09	136	687471	20.00	ng	-0.02
39) Acenaphthene-d10	9.84	164	336741	20.00	ng	-0.02
64) Phenanthrene-d10	11.33	188	529688	20.00	ng	-0.01
76) Chrysene-d12	13.96	240	406889	20.00	ng	-0.02
87) Perylene-d12	15.42	264	510399	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	55472	3.95	ng	0.00
7) Phenol-d6	6.44	99	73797	4.23	ng	-0.01
23) Nitrobenzene-d5	7.36	82	44911	3.89	ng	-0.02
42) 2,4,6-Tribromophenol	10.63	330	12438	3.09	ng	-0.02
45) 2-Fluorobiphenyl	9.16	172	70347	3.73	ng	-0.02
79) Terphenyl-d14	12.90	244	64965	3.28	ng	-0.02
Target Compounds						
31) Naphthalene	8.11	128	2496820	79.410	ng	99
37) 2-Methylnaphthalene	8.80	142	789289	36.195	ng	98
38) 1-Methylnaphthalene	8.90	142	580380	28.292	ng	100
46) 1,1'-Biphenyl	9.26	154	148743	6.385	ng	96
49) Acenaphthylene	9.70	152	303653	10.860	ng	97
52) Acenaphthene	9.87	154	64054	3.477	ng	95
55) Dibenzofuran	10.05	168	167815	6.302	ng	# 93
58) Fluorene	10.39	166	392789	19.608	ng	100
71) Phenanthrene	11.35	178	1659448	63.674	ng	98
72) Anthracene	11.40	178	359470	13.865	ng	96
73) Carbazole	11.55	167	61632	2.709	ng	98
75) Fluoranthene	12.54	202	870925	30.608	ng	99
78) Pvrene	12.77	202	1073183	38.447	ng	98
81) Benzo(a)anthracene	13.95	228	471593	19.477	ng	93
83) Chrysene	13.99	228	443944	18.269	ng	97
86) Indeno(1,2,3-cd)pyrene	16.85	276	269698	11.046	ng	97
88) Benzo(b)fluoranthene	15.00	252	466684m	14.735	ng	
89) Benzo(k)fluoranthene	15.02	252	167675m	6.080	ng	
90) Benzo(a)pvrene	15.35	252	484848	17.888	ng	99
91) Dibenzo(a,h)anthracene	16.85	278	88098m	3.661	ng	
92) Benzo(g,h,i)perylene	17.29	276	294637	12.733	ng	99

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

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