

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012120\
 Data File : BF118660.D
 Acq On : 22 Jan 2020 11:38
 Operator : CG/JU
 Sample : L1148-03DL 100X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-12-(4-6)DL

Manual Integrations
 APPROVED

mohammad
 1/22/2020 1:37:01 PM

Quant Time: Jan 22 12:54:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 20 14:45:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	341666	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	1308404	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	690559	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	1117646	20.00	ng	0.00
76) Chrysene-d12	14.04	240	686654	20.00	ng	0.00
87) Perylene-d12	15.51	264	896201	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.49	112	18588	1.01	ng	-0.01
7) Phenol-d6	6.49	99	26364	1.10	ng	-0.02
23) Nitrobenzene-d5	7.43	82	14212	0.61	ng	-0.02
42) 2,4,6-Tribromophenol	10.70	330	5557	0.70	ng	-0.01
45) 2-Fluorobiphenyl	9.23	172	33545	0.80	ng	-0.01
79) Terphenyl-d14	12.98	244	33156	1.05	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
31) Naphthalene	8.19	128	4807703	81.342	ng	96
37) 2-Methylnaphthalene	8.87	142	976623	23.409	ng	98
38) 1-Methylnaphthalene	8.97	142	592656	14.953	ng	100
46) 1,1'-Biphenyl	9.33	154	263456	5.536	ng	96
49) Acenaphthylene	9.77	152	423540	7.505	ng	96
52) Acenaphthene	9.95	154	486189	12.884	ng	99
55) Dibenzofuran	10.12	168	859283	16.431	ng	92
58) Fluorene	10.46	166	1134856	27.643	ng	99
71) Phenanthrene	11.43	178	3595948	66.690	ng	99
72) Anthracene	11.47	178	1184726	22.205	ng	94
73) Carbazole	11.62	167	356633	7.287	ng	95
75) Fluoranthene	12.62	202	2388743	42.003	ng	98
78) Pvrene	12.85	202	2052718	44.033	ng	94
81) Benzo(a)anthracene	14.03	228	950875m	23.020	ng	
83) Chrysene	14.06	228	777478	19.639	ng	95
86) Indeno(1,2,3-cd)pyrene	16.98	276	472938	13.749	ng	98
88) Benzo(b)fluoranthene	15.08	252	943830m	16.755	ng	
89) Benzo(k)fluoranthene	15.10	252	397420m	7.615	ng	
90) Benzo(a)pvrene	15.44	252	852035	17.534	ng	99
91) Dibenzo(a,h)anthracene	16.99	278	132614	3.085	ng	98
92) Benzo(g,h,i)perylene	17.43	276	428861	10.276	ng	98

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

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