

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\
 Data File : BF121491.D
 Acq On : 31 Aug 2020 22:05
 Operator : JU/CG
 Sample : L3847-03 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LINDEN-JOP-BH-08-B

Manual Integrations
 APPROVED

mohammad
 9/2/2020 1:56:31 PM

Quant Time: Sep 01 09:46:46 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 28 09:25:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	304631	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	1117621	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	527137	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	855498	20.00	ng	0.00
76) Chrysene-d12	14.04	240	607011	20.00	ng	0.01
86) Perylene-d12	15.52	264	628513	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	170225	9.38	ng	0.00
7) Phenol-d6	6.48	99	223451	8.81	ng	-0.01
23) Nitrobenzene-d5	7.42	82	134062	8.20	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	47658	8.95	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	284585	8.71	ng	0.00
79) Terphenyl-d14	12.97	244	222768	7.73	ng	0.00
Target Compounds						
31) Naphthalene	8.17	128	140790	2.537	ng	98
37) 2-Methylnaphthalene	8.86	142	138446	3.766	ng	99
38) 1-Methylnaphthalene	8.96	142	202168	5.824	ng	99
49) Acenaphthylene	9.76	152	201193	4.102	ng	# 94
52) Acenaphthene	9.93	154	329027	10.625	ng	96
58) Fluorene	10.45	166	359254	10.594	ng	99
71) Phenanthrene	11.42	178	3408655	77.642	ng	98
72) Anthracene	11.47	178	1016167	23.533	ng	97
75) Fluoranthene	12.62	202	4586484	96.934	ng	97
78) Pyrene	12.85	202	6177205	142.491	ng	93
81) Benzo(a)anthracene	14.03	228	3037196	81.965	ng	94
83) Chrysene	14.07	228	3188387	86.005	ng	96
87) Indeno(1,2,3-cd)pyrene	17.00	276	1227184	31.734	ng	97
88) Benzo(b)fluoranthene	15.09	252	2991827m	73.177	ng	
89) Benzo(k)fluoranthene	15.12	252	757517m	19.920	ng	
90) Benzo(a)pyrene	15.46	252	2437634	66.306	ng	98
91) Dibenzo(a,h)anthracene	17.00	278	353906	11.180	ng	93
92) Benzo(g,h,i)perylene	17.44	276	1265638	41.505	ng	97

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

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