

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\
 Data File : BF121504.D
 Acq On : 1 Sep 2020 19:08
 Operator : JU/CG
 Sample : L3848-23
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleID :
 LINDEN-BH-01-C

Manual Integrations
 APPROVED

mohammad
 9/2/2020 1:57:23 PM

Quant Time: Sep 02 02:00:23 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	295521	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1124868	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	561459	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	953150	20.00	ng	0.00
76) Chrysene-d12	14.03	240	567467	20.00	ng	0.00
86) Perylene-d12	15.52	264	705697	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1343386	76.27	ng	0.00
7) Phenol-d6	6.48	99	1762265	71.59	ng	-0.01
23) Nitrobenzene-d5	7.42	82	1135917	69.04	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	482569	85.09	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	2050321	58.92	ng	-0.01
79) Terphenyl-d14	12.97	244	1616434	59.96	ng	0.00
Target Compounds						
49) Acenaphthylene	9.76	152	238406	4.564	ng	96
50) Dimethylphthalate	9.61	163	194772	5.050	ng	98
52) Acenaphthene	9.93	154	140288	4.253	ng	98
55) Dibenzofuran	10.10	168	101678	2.122	ng	# 96
58) Fluorene	10.45	166	178734	4.949	ng	100
71) Phenanthrene	11.42	178	2674654	54.681	ng	99
72) Anthracene	11.46	178	784567	16.308	ng	96
73) Carbazole	11.61	167	194235	4.187	ng	98
75) Fluoranthene	12.61	202	4072268	77.248	ng	98
78) Pyrene	12.84	202	4164426	102.756	ng	98
81) Benzo(a)anthracene	14.03	228	2336810	67.458	ng	97
83) Chrysene	14.06	228	2134368	61.585	ng	98
87) Indeno(1,2,3-cd)pyrene	17.00	276	1367116	31.486	ng	98
88) Benzo(b)fluoranthene	15.09	252	2731363m	59.500	ng	
89) Benzo(k)fluoranthene	15.11	252	814793m	19.083	ng	
90) Benzo(a)pyrene	15.46	252	2167180	52.502	ng	96
91) Dibenzo(a,h)anthracene	17.00	278	361370	10.167	ng	92
92) Benzo(g,h,i)perylene	17.44	276	1253125	36.600	ng	97

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

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