

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\
 Data File : BF121484.D
 Acq On : 31 Aug 2020 18:31
 Operator : JU/CG
 Sample : L3847-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LINDEN-JOP-BH-10-A

Manual Integrations
 APPROVED

mohammad
 9/2/2020 1:55:59 PM

Quant Time: Aug 31 19:02:55 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	319306	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	1145636	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	562369	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	947490	20.00	ng	0.00
76) Chrysene-d12	14.03	240	640489	20.00	ng	0.00
86) Perylene-d12	15.50	264	716497	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	1845712	96.98	ng	0.01
7) Phenol-d6	6.49	99	2462210	92.58	ng	0.00
23) Nitrobenzene-d5	7.42	82	1585110	94.59	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	611944	107.73	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	2681305	76.93	ng	0.00
79) Terphenyl-d14	12.97	244	2240719	73.64	ng	0.00
Target Compounds						Qvalue
49) Acenaphthylene	9.76	152	134214	2.565	ng	97
50) Dimethylphthalate	9.61	163	147388	3.815	ng	98
52) Acenaphthene	9.93	154	122118	3.696	ng	99
55) Dibenzofuran	10.10	168	131800	2.746	ng	95
58) Fluorene	10.45	166	191574	5.295	ng	99
71) Phenanthrene	11.41	178	2338148	48.087	ng	100
72) Anthracene	11.46	178	772234	16.148	ng	96
73) Carbazole	11.61	167	185732	4.028	ng	97
75) Fluoranthene	12.60	202	3394296	64.772	ng	99
78) Pyrene	12.83	202	3487897	76.251	ng	98
81) Benzo(a)anthracene	14.02	228	1889742	48.333	ng	97
83) Chrysene	14.06	228	1726604	44.140	ng	98
87) Indeno(1,2,3-cd)pyrene	16.97	276	901574	20.451	ng	96
88) Benzo(b)fluoranthene	15.07	252	2194503m	47.084	ng	
89) Benzo(k)fluoranthene	15.10	252	660979m	15.247	ng	
90) Benzo(a)pyrene	15.44	252	1662736	39.674	ng	96
91) Dibenzo(a,h)anthracene	16.97	278	236690	6.559	ng	94
92) Benzo(g,h,i)perylene	17.41	276	833072	23.965	ng	97

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

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