

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\
 Data File : BF121485.D
 Acq On : 31 Aug 2020 19:02
 Operator : JU/CG
 Sample : L3847-27
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LINDEN-JOP-BH-12-B

Manual Integrations
 APPROVED

mohammad
 9/2/2020 1:56:07 PM

Quant Time: Sep 01 01:04:37 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	285747	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	1044917	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	534528	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	925340	20.00	ng	0.00
76) Chrysene-d12	14.03	240	641348	20.00	ng	0.00
86) Perylene-d12	15.51	264	668094	20.00	ng	0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1067907	62.70	ng	0.00
7) Phenol-d6	6.48	99	1382231	58.07	ng	-0.01
23) Nitrobenzene-d5	7.42	82	894237	58.51	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	368883	68.32	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1632013	49.26	ng	0.00
79) Terphenyl-d14	12.97	244	1343101	44.08	ng	0.00

Target Compounds				Qvalue		
49) Acenaphthylene	9.76	152	123046	2.474	ng	98
50) Dimethylphthalate	9.61	163	125233	3.411	ng	100
52) Acenaphthene	9.93	154	101500	3.232	ng	98
58) Fluorene	10.45	166	118387	3.443	ng	99
71) Phenanthrene	11.41	178	1846442	38.883	ng	99
72) Anthracene	11.46	178	513669	10.998	ng	96
73) Carbazole	11.61	167	175946	3.907	ng	99
75) Fluoranthene	12.60	202	2704813	52.851	ng	100
78) Pyrene	12.83	202	2481640	54.180	ng	99
81) Benzo(a)anthracene	14.02	228	1374927	35.118	ng	97
83) Chrysene	14.06	228	1110910	28.362	ng	96
87) Indeno(1,2,3-cd)pyrene	16.99	276	660428	16.066	ng	96
88) Benzo(b)fluoranthene	15.08	252	1544808m	35.546	ng	
89) Benzo(k)fluoranthene	15.10	252	440214m	10.890	ng	
90) Benzo(a)pyrene	15.45	252	1095512	28.033	ng	97
91) Dibenzo(a,h)anthracene	16.99	278	167733	4.985	ng	# 88
92) Benzo(g,h,i)perylene	17.43	276	669891	20.667	ng	97

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

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