

Data File : BF121510.D
Acq On : 1 Sep 2020 22:15
Operator : JU/CG
Sample : L3847-03DL 25X
Misc :
ALS Vial : 17 Sample Multiplier: 1

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

Manual Integrations
APPROVED

mohammad
9/2/2020 1:57:39 PM

Quant Time: Sep 07 04:10:23 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	262532	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1020775	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	518443	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	887889	20.00	ng	0.00
76) Chrysene-d12	14.03	240	644269	20.00	ng	0.00
86) Perylene-d12	15.51	264	714210	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	24969	1.60	ng	-0.01
7) Phenol-d6	6.47	99	36741	1.68	ng	-0.02
23) Nitrobenzene-d5	7.42	82	20811	1.39	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	9135	1.74	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	55316	1.72	ng	-0.01
79) Terphenyl-d14	12.97	244	49553	1.62	ng	0.00
Target Compounds						Qvalue
52) Acenaphthene	9.93	154	69065	2.268	ng	98
71) Phenanthrene	11.41	178	960270	21.075	ng	98
72) Anthracene	11.46	178	224419	5.008	ng	95
75) Fluoranthene	12.60	202	1376112	28.023	ng	97
78) Pyrene	12.83	202	1896306	41.213	ng	98
81) Benzo(a)anthracene	14.02	228	691151	17.573	ng	98
83) Chrysene	14.06	228	784388m	19.935	ng	
87) Indeno(1,2,3-cd)pyrene	16.97	276	309793	7.050	ng	95
88) Benzo(b)fluoranthene	15.08	252	675425m	14.538	ng	
89) Benzo(k)fluoranthene	15.09	252	235093m	5.440	ng	
90) Benzo(a)pyrene	15.44	252	628796	15.052	ng	99
91) Dibenzo(a,h)anthracene	16.99	278	86932m	2.417	ng	
92) Benzo(g,h,i)perylene	17.42	276	306939	8.858	ng	98

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

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