

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031117\
 Data File : BM009205.D
 Acq On : 11 Mar 2017 13:31
 Operator : SJ/MA
 Sample : I2107-07DL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AJ6DL

Manual Integrations
 APPROVED

mohammad
 3/13/2017 5:13:22 PM

Quant Time: Mar 13 11:37:44 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 13 11:21:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	257	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	861	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	525	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	1142	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1287	0.40	ng/ul	0.00
20) Perylene-d12	23.77	264	1236	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	30	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	67	0.02	ng/ul	0.00
Target Compounds						
						Qvalue
3) Naphthalene	10.73	128	2057	0.83	ng/ul	99
5) 2-Methylnaphthalene	12.34	142	2085	1.17	ng/ul	99
7) Acenaphthylene	14.23	152	343	0.13	ng/ul#	54
8) Acenaphthene	14.57	153	95	0.05	ng/ul#	81
9) Fluorene	15.56	166	139	0.06	ng/ul#	56
12) Phenanthrene	17.30	178	2787	0.78	ng/ul	99
13) Anthracene	17.40	178	478m	0.13	ng/ul	
15) Fluoranthene	19.33	202	4751	1.02	ng/ul	93
17) Pyrene	19.69	202	4455	1.01	ng/ul	95
18) Benzo(a)anthracene	21.42	228	2733	0.64	ng/ul	95
19) Chrysene	21.47	228	3482	0.86	ng/ul	94
21) Benzo(b)fluoranthene	23.07	252	5957m	1.23	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	1399m	0.33	ng/ul	
23) Benzo(a)pyrene	23.67	252	2945	0.67	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.15	276	2433	0.44	ng/ul#	88
25) Dibenzo(a,h)anthracene	26.16	278	612	0.13	ng/ul#	65
26) Benzo(g,h,i)perylene	26.88	276	2187	0.46	ng/ul#	91

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

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