

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030917\
 Data File : BM009161.D
 Acq On : 10 Mar 2017 00:33
 Operator : SJ/MA
 Sample : I2062-21DL 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AK7DL

Manual Integrations
 APPROVED

mohammad
 3/10/2017 6:01:48 PM

Quant Time: Mar 10 04:02:03 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 10 03:16:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	263	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	798	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	441	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	1021	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1142	0.40	ng/ul	0.00
20) Perylene-d12	23.77	264	1170	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	75	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	199	0.06	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	385	0.17	ng/ul#	71
5) 2-Methylnaphthalene	12.34	142	282	0.17	ng/ul	95
7) Acenaphthylene	14.23	152	180	0.08	ng/ul#	34
12) Phenanthrene	17.30	178	1180	0.37	ng/ul#	90
13) Anthracene	17.40	178	385	0.12	ng/ul#	73
15) Fluoranthene	19.33	202	3315	0.79	ng/ul	94
17) Pyrene	19.69	202	3311	0.84	ng/ul	98
18) Benzo(a)anthracene	21.42	228	2739m	0.72	ng/ul	
19) Chrysene	21.47	228	3161m	0.88	ng/ul	
21) Benzo(b)fluoranthene	23.07	252	5673m	1.23	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	1948m	0.48	ng/ul	
23) Benzo(a)pyrene	23.67	252	3089	0.75	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.15	276	2194	0.42	ng/ul#	88
25) Dibenzo(a,h)anthracene	26.16	278	584	0.13	ng/ul#	63
26) Benzo(g,h,i)perylene	26.89	276	1902	0.42	ng/ul#	92

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

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