

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031017\
 Data File : BM009174.D
 Acq On : 10 Mar 2017 13:58
 Operator : SJ/MA
 Sample : I2062-22
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AK8

Manual Integrations
 APPROVED

mohammad
 3/13/2017 2:13:28 PM

Quant Time: Mar 10 16:54:13 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 15:19:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	264	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	830	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	491	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	1098m	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1194	0.40	ng/ul	0.00
20) Perylene-d12	23.77	264	1028m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	277	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	769	0.22	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	1315	0.55	ng/ul	96
5) 2-Methylnaphthalene	12.34	142	1256	0.73	ng/ul	100
7) Acenaphthylene	14.23	152	253	0.10	ng/ul#	41
8) Acenaphthene	14.57	153	70	0.04	ng/ul#	79
12) Phenanthrene	17.30	178	2444	0.71	ng/ul	98
13) Anthracene	17.40	178	262	0.08	ng/ul#	63
15) Fluoranthene	19.33	202	5804	1.29	ng/ul	94
17) Pyrene	19.69	202	5036	1.23	ng/ul	95
18) Benzo(a)anthracene	21.42	228	2557	0.64	ng/ul	94
19) Chrysene	21.47	228	3600m	0.96	ng/ul	
21) Benzo(b)fluoranthene	23.07	252	5627m	1.39	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	1394m	0.39	ng/ul	
23) Benzo(a)pyrene	23.67	252	3276m	0.90	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.15	276	2449m	0.53	ng/ul	
25) Dibenzo(a,h)anthracene	26.16	278	640m	0.16	ng/ul	
26) Benzo(g,h,i)perylene	26.89	276	2278m	0.58	ng/ul	

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

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