

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030917\
 Data File : BM009144.D
 Acq On : 09 Mar 2017 14:21
 Operator : SJ/MA
 Sample : I2062-20
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AK6

Manual Integrations
 APPROVED

mohammad
 3/10/2017 6:00:56 PM

Quant Time: Mar 09 15:39:34 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 09 15:10:54 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	194	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	630	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.52	164	392	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	889m	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1111	0.40	ng/ul	-0.01
20) Perylene-d12	23.78	264	914m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	285	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	904	0.33	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	4582	2.52	ng/ul	97
5) 2-Methylnaphthalene	12.34	142	3768	2.88	ng/ul	99
7) Acenaphthylene	14.23	152	359	0.18	ng/ul#	44
8) Acenaphthene	14.59	153	146	0.09	ng/ul#	79
9) Fluorene	15.56	166	139	0.08	ng/ul#	41
12) Phenanthrene	17.31	178	4475	1.60	ng/ul	92
13) Anthracene	17.40	178	664	0.24	ng/ul#	86
15) Fluoranthene	19.33	202	8666	2.38	ng/ul	94
17) Pyrene	19.69	202	9359	2.45	ng/ul	97
18) Benzo(a)anthracene	21.42	228	6333	1.72	ng/ul	95
19) Chrysene	21.48	228	5872	1.68	ng/ul	95
21) Benzo(b)fluoranthene	23.07	252	10188m	2.84	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	3255m	1.03	ng/ul	
23) Benzo(a)pyrene	23.67	252	5661m	1.75	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.15	276	4133m	1.01	ng/ul	
25) Dibenzo(a,h)anthracene	26.16	278	1157m	0.34	ng/ul	
26) Benzo(g,h,i)perylene	26.89	276	3639m	1.04	ng/ul	

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

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