

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030817\
 Data File : BM009125.D
 Acq On : 09 Mar 2017 02:48
 Operator : SJ/MA
 Sample : I2062-20
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AK6

Manual Integrations
 APPROVED

mohammad
 3/9/2017 2:01:53 PM

Quant Time: Mar 09 05:41:18 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 02:44:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	358	0.40	ng/ul	0.02
2) Naphthalene-d8	10.69	136	1248	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	805	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.28	188	2058m	0.40	ng/ul	0.02
16) Chrysene-d12	21.45	240	2354	0.40	ng/ul	0.00
20) Perylene-d12	23.80	264	1659m	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	531	0.26	ng/ul	0.01
14) Fluoranthene-d10	19.30	212	1857	0.29	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	8261	2.30	ng/ul	95
5) 2-Methylnaphthalene	12.35	142	7108	2.75	ng/ul	100
7) Acenaphthylene	14.25	152	631	0.15	ng/ul#	66
8) Acenaphthene	14.59	153	289	0.09	ng/ul#	85
9) Fluorene	15.58	166	289	0.08	ng/ul#	82
12) Phenanthrene	17.31	178	9350	1.45	ng/ul	96
13) Anthracene	17.40	178	1313	0.21	ng/ul#	93
15) Fluoranthene	19.33	202	17996	2.14	ng/ul	93
17) Pyrene	19.69	202	18586	2.30	ng/ul	97
18) Benzo(a)anthracene	21.43	228	11039	1.41	ng/ul	100
19) Chrysene	21.48	228	11658	1.57	ng/ul	97
21) Benzo(b)fluoranthene	23.08	252	17530m	2.69	ng/ul	
22) Benzo(k)fluoranthene	23.12	252	4637m	0.80	ng/ul	
23) Benzo(a)pyrene	23.68	252	9724m	1.66	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.17	276	5713m	0.77	ng/ul	
25) Dibenzo(a,h)anthracene	26.18	278	1449m	0.23	ng/ul	
26) Benzo(g,h,i)perylene	26.91	276	5064m	0.80	ng/ul	

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

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