

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030817\
 Data File : BM009114.D
 Acq On : 08 Mar 2017 18:05
 Operator : SJ/MA
 Sample : I2062-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC8

Manual Integrations
 APPROVED

mohammad
 3/9/2017 2:01:18 PM

Quant Time: Mar 09 01:36:48 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 01:19:41 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	217	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.69	136	668	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	389	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	981m	0.40	ng/ul	0.00
16) Chrysene-d12	21.45	240	977	0.40	ng/ul	0.00
20) Perylene-d12	23.78	264	879m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	304	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	805	0.26	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.31	178	604m	0.20	ng/ul	
15) Fluoranthene	19.33	202	1424	0.35	ng/ul	95
17) Pyrene	19.69	202	1563	0.47	ng/ul	96
18) Benzo(a)anthracene	21.43	228	1744	0.54	ng/ul	96
19) Chrysene	21.48	228	2056	0.67	ng/ul	97
21) Benzo(b)fluoranthene	23.07	252	3184m	0.92	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	1258m	0.41	ng/ul	
23) Benzo(a)pyrene	23.67	252	2400m	0.77	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.16	276	1323m	0.33	ng/ul	
25) Dibenzo(a,h)anthracene	26.17	278	377m	0.11	ng/ul	
26) Benzo(g,h,i)perylene	26.90	276	1146m	0.34	ng/ul	

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

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