

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021215\
 Data File : BM000508.D
 Acq On : 12 Feb 2015 18:27
 Operator : TP/IZ
 Sample : G1249-23
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4FN6

Manual Integrations
 APPROVED

mohammad
 2/13/2015 5:41:21 PM

Quant Time: Feb 13 10:36:07 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 12 16:45:55 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	3562	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	12052	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	7275	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	14903m	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	12385	0.40	ng/ul	-0.01
20) Perylene-d12	23.51	264	10277m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	6839	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	15045m	0.48	ng/ul	0.00
Target Compounds				Qvalue		
8) Acenaphthene	14.41	153	2438	0.12	ng/ul	97
9) Fluorene	15.40	166	1173	0.05	ng/ul#	95
11) Pentachlorophenol	16.75	266	316	0.17	ng/ul	98
12) Phenanthrene	17.13	178	2742m	0.07	ng/ul	
13) Anthracene	17.23	178	3134	0.09	ng/ul#	85
15) Fluoranthene	19.15	202	19200	0.43	ng/ul	92
17) Pyrene	19.53	202	23928	0.54	ng/ul	98
18) Benzo(a)anthracene	21.27	228	2238	0.06	ng/ul#	87
19) Chrysene	21.32	228	2153	0.05	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.75	276	1290	0.04	ng/ul	97
25) Dibenzo(a,h)anthracene	25.77	278	1006m	0.04	ng/ul	
26) Benzo(g,h,i)perylene	26.44	276	1184	0.04	ng/ul#	74

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

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