

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021115\
 Data File : BM000488.D
 Acq On : 11 Feb 2015 22:10
 Operator : TP/IZ
 Sample : G1249-04MSD
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4FL2MSD

Manual Integrations
 APPROVED

apatel
 2/12/2015 4:10:25 PM

Quant Time: Feb 12 15:05:05 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 11 23:58:48 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	3186	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	11446	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	7546	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	15792m	0.40	ng/ul	0.00
16) Chrysene-d12	21.30	240	11934	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	9307m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	6220	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.14	212	14305m	0.41	ng/ul	0.00
Target Compounds						
8) Acenaphthene	14.41	153	7710	0.33	ng/ul	97
11) Pentachlorophenol	16.76	266	1199	0.82	ng/ul	98
12) Phenanthrene	17.15	178	304m	0.01	ng/ul	
15) Fluoranthene	19.16	202	833	0.02	ng/ul	82
17) Pyrene	19.53	202	23108	0.53	ng/ul#	93
18) Benzo(a)anthracene	21.28	228	1292	0.04	ng/ul#	89
19) Chrysene	21.33	228	1406	0.03	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.78	276	843	0.03	ng/ul	96
26) Benzo(g,h,i)perylene	26.47	276	756	0.03	ng/ul#	62

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

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