

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021215\
 Data File : BM000512.D
 Acq On : 12 Feb 2015 20:49
 Operator : TP/IZ
 Sample : G1249-04MSD
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4FL2MSD

Manual Integrations
 APPROVED

mohammad
 2/13/2015 5:41:32 PM

Quant Time: Feb 13 10:59:42 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	3781	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	12781	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	8176	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	15827m	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	12850	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	10867m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	6550	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	14325m	0.43	ng/ul	0.00
Target Compounds						
8) Acenaphthene	14.41	153	7900	0.33	ng/ul	97
11) Pentachlorophenol	16.75	266	1250	0.62	ng/ul	96
15) Fluoranthene	19.16	202	879	0.02	ng/ul	72
17) Pyrene	19.53	202	22912	0.50	ng/ul	99
18) Benzo(a)anthracene	21.27	228	1410	0.04	ng/ul#	91
19) Chrysene	21.32	228	1518	0.04	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.75	276	1255	0.03	ng/ul	98
25) Dibenzo(a,h)anthracene	25.77	278	1001m	0.03	ng/ul	
26) Benzo(g,h,i)perylene	26.44	276	1146	0.03	ng/ul#	71

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

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