

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021115\
 Data File : BM000487.D
 Acq On : 11 Feb 2015 21:34
 Operator : TP/IZ
 Sample : G1249-03MS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4FL2MS

Manual Integrations
 APPROVED

apatel
 2/12/2015 4:10:21 PM

Quant Time: Feb 12 15:03:38 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	3292	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	11438	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	7410	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	14417m	0.40	ng/ul	0.00
16) Chrysene-d12	21.30	240	10135	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	7392m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	6832	0.45	ng/ul	0.00
14) Fluoranthene-d10	19.14	212	14201m	0.44	ng/ul	0.00
Target Compounds						Qvalue
8) Acenaphthene	14.41	153	7625	0.34	ng/ul	95
11) Pentachlorophenol	16.76	266	1078	0.81	ng/ul	99
12) Phenanthrene	17.15	178	331m	0.01	ng/ul	
15) Fluoranthene	19.16	202	841	0.02	ng/ul	77
17) Pyrene	19.53	202	20818	0.56	ng/ul#	94
18) Benzo(a)anthracene	21.28	228	1234	0.04	ng/ul#	90
19) Chrysene	21.33	228	1361	0.04	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.77	276	874	0.03	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.80	278	716	0.04	ng/ul#	32
26) Benzo(g,h,i)perylene	26.46	276	826	0.04	ng/ul#	66

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

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