

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090315\
 Data File : BM002651.D
 Acq On : 03 Sep 2015 17:00
 Operator : UM/IZ
 Sample : MDL-01-W
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-01-W

Quant Time: Sep 04 04:22:13 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM90315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 03 14:33:37 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.78	152	7597	0.40	ng/ul	0.00
4) Naphthalene-d8	11.64	136	32111	0.40	ng/ul	0.00
8) Acenaphthene-d10	15.36	164	14548	0.40	ng/ul	0.00
12) Phenanthrene-d10	18.08	188	26099	0.40	ng/ul	0.00
18) Chrysene-d12	22.17	240	19261	0.40	ng/ul	0.01
22) Perylene-d12	24.78	264	18967	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
6) 2-Methylnaphthalene-d10	13.17	152	17068	0.40	ng/ul	0.00
16) Fluoranthene-d10	20.05	212	22783	0.38	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
5) Naphthalene	11.69	128	4187	0.05	ng/ul	96
7) 2-Methylnaphthalene	13.25	142	3371	0.06	ng/ul	94
9) Acenaphthylene	15.09	152	3908	0.05	ng/ul	97
10) Acenaphthene	15.42	153	2769	0.05	ng/ul	99
11) Fluorene	16.39	166	2972	0.05	ng/ul	94
14) Phenanthrene	18.11	178	4119	0.05	ng/ul	94
15) Anthracene	18.20	178	3935	0.05	ng/ul	95
17) Fluoranthene	20.07	202	4217	0.05	ng/ul	91
19) Pyrene	20.42	202	4205	0.06	ng/ul#	86
20) Benzo(a)anthracene	22.14	228	3626	0.06	ng/ul#	82
21) Chrysene	22.20	228	3588	0.06	ng/ul#	87
23) Benzo(b)fluoranthene	23.97	252	4087	0.06	ng/ul#	42
24) Benzo(k)fluoranthene	24.02	252	3320	0.05	ng/ul#	39
25) Benzo(a)pyrene	24.67	252	3449	0.06	ng/ul#	37
26) Indeno(1,2,3-cd)pyrene	27.56	276	4108	0.06	ng/ul	98
27) Dibenzo(a,h)anthracene	27.56	278	3395	0.06	ng/ul#	57
28) Benzo(g,h,i)perylene	28.42	276	3516	0.06	ng/ul#	66

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

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