

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090315\
 Data File : BM002652.D
 Acq On : 03 Sep 2015 17:37
 Operator : UM/IZ
 Sample : MDL-02-W
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-02-W

Quant Time: Sep 04 04:23:05 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	7551	0.40	ng/ul	0.00
4) Naphthalene-d8	11.64	136	29130	0.40	ng/ul	0.00
8) Acenaphthene-d10	15.36	164	11726	0.40	ng/ul	0.00
12) Phenanthrene-d10	18.08	188	20761	0.40	ng/ul	0.00
18) Chrysene-d12	22.17	240	19201	0.40	ng/ul	0.01
22) Perylene-d12	24.78	264	18905	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
6) 2-Methylnaphthalene-d10	13.17	152	17546	0.45	ng/ul	0.00
16) Fluoranthene-d10	20.05	212	23853	0.50	ng/ul	0.00
Target Compounds				Qvalue		
5) Naphthalene	11.69	128	4902	0.06	ng/ul	95
7) 2-Methylnaphthalene	13.25	142	3653	0.07	ng/ul	92
9) Acenaphthylene	15.09	152	4305	0.07	ng/ul	99
10) Acenaphthene	15.42	153	2812	0.06	ng/ul	99
11) Fluorene	16.39	166	2986	0.06	ng/ul#	94
14) Phenanthrene	18.11	178	4238	0.07	ng/ul	94
15) Anthracene	18.20	178	4075	0.07	ng/ul	96
17) Fluoranthene	20.07	202	4703	0.07	ng/ul	92
19) Pyrene	20.42	202	4772	0.07	ng/ul#	89
20) Benzo(a)anthracene	22.14	228	4441	0.07	ng/ul#	88
21) Chrysene	22.20	228	4523	0.07	ng/ul#	91
23) Benzo(b)fluoranthene	23.97	252	4671	0.07	ng/ul#	56
24) Benzo(k)fluoranthene	24.02	252	4559	0.07	ng/ul#	54
25) Benzo(a)pyrene	24.67	252	4334	0.07	ng/ul#	50
26) Indeno(1,2,3-cd)pyrene	27.55	276	5230	0.07	ng/ul	96
27) Dibenzo(a,h)anthracene	27.56	278	4076	0.07	ng/ul#	69
28) Benzo(g,h,i)perylene	28.41	276	4366	0.07	ng/ul#	74

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

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