

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090315\
 Data File : BM002660.D
 Acq On : 03 Sep 2015 22:33
 Operator : UM/IZ
 Sample : MDL-01-S
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-01-S

Quant Time: Sep 04 04:41:16 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 : Fri Sep 04 04:35:12 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.78	152	7280	0.40	ng/ul	0.00
4) Naphthalene-d8	11.64	136	29075	0.40	ng/ul	0.00
8) Acenaphthene-d10	15.36	164	12178	0.40	ng/ul	0.00
12) Phenanthrene-d10	18.08	188	21167	0.40	ng/ul	0.00
18) Chrysene-d12	22.16	240	18319	0.40	ng/ul	0.00
22) Perylene-d12	24.77	264	18030	0.40	ng/ul	-0.01

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	17395	0.45	ng/ul	0.00
16) Fluoranthene-d10	20.03	212	23258	0.48	ng/ul	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
5) Naphthalene	11.69	128	4522	0.06	ng/ul	95
7) 2-Methylnaphthalene	13.25	142	3258	0.06	ng/ul	95
9) Acenaphthylene	15.09	152	4036	0.06	ng/ul	98
10) Acenaphthene	15.42	153	2748	0.06	ng/ul	98
11) Fluorene	16.39	166	2924	0.06	ng/ul#	95
14) Phenanthrene	18.11	178	4025	0.06	ng/ul	95
15) Anthracene	18.20	178	3778	0.06	ng/ul	95
17) Fluoranthene	20.07	202	4383	0.07	ng/ul	89
19) Pyrene	20.42	202	4488	0.06	ng/ul#	93
20) Benzo(a)anthracene	22.14	228	4082	0.07	ng/ul#	91
21) Chrysene	22.19	228	3898	0.07	ng/ul#	89
23) Benzo(b)fluoranthene	23.96	252	4622	0.07	ng/ul#	56
24) Benzo(k)fluoranthene	24.01	252	3792	0.06	ng/ul#	51
25) Benzo(a)pyrene	24.66	252	3797	0.06	ng/ul#	47
26) Indeno(1,2,3-cd)pyrene	27.54	276	4545	0.06	ng/ul	97
27) Dibenzo(a,h)anthracene	27.55	278	3692	0.06	ng/ul#	68
28) Benzo(g,h,i)perylene	28.40	276	3934	0.07	ng/ul#	75

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

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