

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090315\
 Data File : BM002654.D
 Acq On : 03 Sep 2015 18:51
 Operator : UM/IZ
 Sample : MDL-04-W
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-04-W

Quant Time: Sep 04 04:24:39 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	7929	0.40	ng/ul	0.00
4) Naphthalene-d8	11.64	136	31158	0.40	ng/ul	0.00
8) Acenaphthene-d10	15.36	164	13067	0.40	ng/ul	0.00
12) Phenanthrene-d10	18.08	188	23490	0.40	ng/ul	0.00
18) Chrysene-d12	22.16	240	22295	0.40	ng/ul	0.00
22) Perylene-d12	24.79	264	21757	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	17957	0.43	ng/ul	0.00
16) Fluoranthene-d10	20.05	212	26735	0.50	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
5) Naphthalene	11.69	128	4198	0.05	ng/ul	94
7) 2-Methylnaphthalene	13.25	142	3236	0.06	ng/ul	88
9) Acenaphthylene	15.09	152	3800	0.05	ng/ul	98
10) Acenaphthene	15.42	153	2565	0.05	ng/ul	99
11) Fluorene	16.39	166	2726	0.05	ng/ul	96
14) Phenanthrene	18.11	178	3874	0.06	ng/ul#	92
15) Anthracene	18.20	178	3778	0.05	ng/ul#	93
17) Fluoranthene	20.07	202	4510	0.06	ng/ul	89
19) Pyrene	20.42	202	4677	0.06	ng/ul#	92
20) Benzo(a)anthracene	22.14	228	4417	0.06	ng/ul#	88
21) Chrysene	22.20	228	4430	0.06	ng/ul#	90
23) Benzo(b)fluoranthene	23.97	252	4814	0.06	ng/ul#	50
24) Benzo(k)fluoranthene	24.02	252	4454	0.06	ng/ul#	46
25) Benzo(a)pyrene	24.67	252	4273	0.06	ng/ul#	43
26) Indeno(1,2,3-cd)pyrene	27.55	276	5033	0.06	ng/ul	98
27) Dibenzo(a,h)anthracene	27.56	278	4004	0.06	ng/ul#	66
28) Benzo(g,h,i)perylene	28.41	276	4322	0.06	ng/ul#	73

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

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