

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM083016\
 Data File : BM007346.D
 Acq On : 30 Aug 2016 17:29
 Operator : UM/SJ
 Sample : MDL-04-W
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-04-W

Quant Time: Aug 30 18:06:53 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM081516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 23:46:46 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	1314	0.40	ng/ul	0.00
2) Naphthalene-d8	10.58	136	5750	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.43	164	3947	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.18	188	10588	0.40	ng/ul	0.00
16) Chrysene-d12	21.35	240	11158	0.40	ng/ul	0.00
20) Perylene-d12	23.62	264	10707	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.18	152	2067	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.20	212	6735	0.28	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.63	128	830	0.06	ng/ul#	72
5) 2-Methylnaphthalene	12.26	142	573	0.06	ng/ul	95
7) Acenaphthylene	14.14	152	1003	0.05	ng/ul#	87
8) Acenaphthene	14.49	153	744	0.05	ng/ul#	88
9) Fluorene	15.48	166	896	0.06	ng/ul#	80
11) Pentachlorophenol	16.83	266	129	0.05	ng/ul	93
12) Phenanthrene	17.21	178	1639	0.05	ng/ul#	89
13) Anthracene	17.30	178	1445	0.05	ng/ul#	85
15) Fluoranthene	19.22	202	2029	0.06	ng/ul	87
17) Pyrene	19.59	202	2161	0.06	ng/ul	93
18) Benzo(a)anthracene	21.33	228	1979	0.06	ng/ul#	90
19) Chrysene	21.38	228	1996	0.06	ng/ul#	90
21) Benzo(b)fluoranthene	22.93	252	2028	0.05	ng/ul	81
22) Benzo(k)fluoranthene	22.98	252	2023	0.06	ng/ul#	81
23) Benzo(a)pyrene	23.52	252	1914	0.06	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.90	276	1990	0.05	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.93	278	1582	0.05	ng/ul#	78
26) Benzo(g,h,i)perylene	26.61	276	1703	0.05	ng/ul#	87

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

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