

Data Path : Z:\HPCHEM1\BNA_N\DATA\BN030518\
 Data File : BN000251.D
 Acq On : 05 Mar 2018 18:35
 Operator : JU/SJ
 Sample : MDL-SIM-S-06
 Misc : 0.07 PPM
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-SIM-S-06

Quant Time: Mar 06 04:38:44 2018
 Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN030218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 06 04:37:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.45	152	4285	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	13486	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.06	164	5697	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.78	188	12900	0.40	ng/ul	0.00
16) Chrysene-d12	21.89	240	7149	0.40	ng/ul	0.00
20) Perylene-d12	24.37	264	5681	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	6699	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.77	212	11316	0.38	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.34	128	2798	0.065	ng/ul#	91
5) 2-Methylnaphthalene	12.92	142	1714	0.063	ng/ul	99
7) Acenaphthylene	14.78	152	1704	0.060	ng/ul	96
8) Acenaphthene	15.12	153	1752	0.064	ng/ul	94
9) Fluorene	16.10	166	1831	0.062	ng/ul	91
11) Pentachlorophenol	17.43	266	97	0.037	ng/ul#	81
12) Phenanthrene	17.82	178	3255	0.062	ng/ul#	90
13) Anthracene	17.91	178	2164	0.055	ng/ul#	91
15) Fluoranthene	19.80	202	3123	0.061	ng/ul	62
17) Pyrene	20.16	202	3041	0.072	ng/ul#	55
18) Benzo(a)anthracene	21.88	228	1808	0.064	ng/ul#	84
19) Chrysene	21.93	228	2701	0.071	ng/ul	95
21) Benzo(b)fluoranthene	23.61	252	2187	0.064	ng/ul#	80
22) Benzo(k)fluoranthene	23.66	252	1916	0.059	ng/ul#	72
23) Benzo(a)pyrene	24.26	252	1593	0.060	ng/ul#	67
24) Indeno(1,2,3-cd)pyrene	26.91	276	1752	0.058	ng/ul#	78
25) Dibenzo(a,h)anthracene	26.93	278	1231	0.057	ng/ul#	59
26) Benzo(g,h,i)perylene	27.70	276	1899	0.064	ng/ul#	71

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

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