

Data Path : Z:\HPCHEM1\BNA_N\DATA\BN030518\
 Data File : BN000239.D
 Acq On : 05 Mar 2018 11:41
 Operator : JU/SJ
 Sample : MDL-SIM-W-03
 Misc : 0.07 PPM
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-SIM-W-03

Quant Time: Mar 05 14:39:37 2018
 Quant Method : Z:\HPCHEM1\BNA_N\Methods\SOM-EPA-SIM-BN030218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 05 14:32:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.45	152	3899	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	12772	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.06	164	5431	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.78	188	12251	0.40	ng/ul	0.00
16) Chrysene-d12	21.89	240	6769	0.40	ng/ul	0.00
20) Perylene-d12	24.37	264	5551	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	6259	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.77	212	10634	0.37	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.34	128	2412	0.059	ng/ul#	91
5) 2-Methylnaphthalene	12.92	142	1491	0.057	ng/ul	99
7) Acenaphthylene	14.79	152	1522	0.056	ng/ul	95
8) Acenaphthene	15.12	153	1479	0.057	ng/ul	95
9) Fluorene	16.10	166	1552	0.055	ng/ul	92
11) Pentachlorophenol	17.43	266	91	0.037	ng/ul	92
12) Phenanthrene	17.82	178	2857	0.057	ng/ul#	88
13) Anthracene	17.91	178	1967	0.053	ng/ul#	91
15) Fluoranthene	19.80	202	2669	0.055	ng/ul	62
17) Pyrene	20.15	202	2588	0.065	ng/ul#	55
18) Benzo(a)anthracene	21.88	228	1487	0.056	ng/ul#	85
19) Chrysene	21.93	228	2431	0.067	ng/ul	95
21) Benzo(b)fluoranthene	23.61	252	1937	0.058	ng/ul#	77
22) Benzo(k)fluoranthene	23.66	252	1759	0.056	ng/ul#	67
23) Benzo(a)pyrene	24.26	252	1478	0.057	ng/ul#	61
24) Indeno(1,2,3-cd)pyrene	26.92	276	1552	0.053	ng/ul#	77
25) Dibenzo(a,h)anthracene	26.93	278	1044	0.049	ng/ul#	59
26) Benzo(g,h,i)perylene	27.71	276	1660	0.057	ng/ul#	71

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

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