

Data Path : Z:\HPCHEM1\BNA_N\DATA\BN030518\
 Data File : BN000237.D
 Acq On : 05 Mar 2018 10:32
 Operator : JU/SJ
 Sample : MDL-SIM-W-01
 Misc : 0.07 PPM
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-SIM-W-01

Quant Time: Mar 05 14:38:24 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 14:31:56 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	2897	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	8780	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.06	164	3802	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.78	188	11351	0.40	ng/ul	0.00
16) Chrysene-d12	21.89	240	7511	0.40	ng/ul	0.00
20) Perylene-d12	24.36	264	6152	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	4185	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.77	212	10451	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.34	128	1728	0.061	ng/ul#	87
5) 2-Methylnaphthalene	12.92	142	1057	0.059	ng/ul	100
7) Acenaphthylene	14.79	152	1165	0.062	ng/ul#	92
8) Acenaphthene	15.12	153	1094	0.060	ng/ul#	93
9) Fluorene	16.10	166	1281	0.065	ng/ul	91
11) Pentachlorophenol	17.42	266	97	0.043	ng/ul	91
12) Phenanthrene	17.82	178	2718	0.059	ng/ul#	89
13) Anthracene	17.91	178	1928	0.056	ng/ul#	91
15) Fluoranthene	19.80	202	2964	0.066	ng/ul	62
17) Pyrene	20.16	202	2880	0.065	ng/ul#	55
18) Benzo(a)anthracene	21.88	228	1766	0.060	ng/ul#	85
19) Chrysene	21.93	228	2758	0.069	ng/ul	94
21) Benzo(b)fluoranthene	23.61	252	2212	0.060	ng/ul	81
22) Benzo(k)fluoranthene	23.66	252	2014	0.058	ng/ul#	70
23) Benzo(a)pyrene	24.26	252	1639	0.057	ng/ul#	65
24) Indeno(1,2,3-cd)pyrene	26.91	276	1831	0.056	ng/ul#	78
25) Dibenzo(a,h)anthracene	26.93	278	1221	0.052	ng/ul#	63
26) Benzo(g,h,i)perylene	27.70	276	2017	0.063	ng/ul#	72

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

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