

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN021419\
 Data File : BN004419.D
 Acq On : 14 Feb 2019 14:28
 Operator : JU/SJ
 Sample : MDL-W-04
 Misc : 0.07 PPM
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-W-04

Quant Time: Feb 14 15:07:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 14:04:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	1196	0.40	ng/ul	0.00
2) Naphthalene-d8	10.53	136	4608	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	2716	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.13	188	7529	0.40	ng/ul	0.00
16) Chrysene-d12	21.33	240	9482	0.40	ng/ul	0.00
20) Perylene-d12	23.60	264	9750	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	2455	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.16	212	7872	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.58	128	823	0.062	ng/ul#	83
5) 2-Methylnaphthalene	12.20	142	590	0.062	ng/ul	100
7) Acenaphthylene	14.11	152	709	0.057	ng/ul#	86
8) Acenaphthene	14.45	153	667	0.062	ng/ul	96
9) Fluorene	15.43	166	795	0.062	ng/ul#	94
11) Pentachlorophenol	16.78	266	58	0.036	ng/ul	88
12) Phenanthrene	17.17	178	1515	0.062	ng/ul#	90
13) Anthracene	17.26	178	1163	0.056	ng/ul#	89
15) Fluoranthene	19.20	202	1882	0.063	ng/ul	94
17) Pyrene	19.56	202	1873	0.061	ng/ul#	91
18) Benzo(a)anthracene	21.31	228	1846	0.060	ng/ul	95
19) Chrysene	21.36	228	2276	0.064	ng/ul	96
21) Benzo(b)fluoranthene	22.91	252	2450	0.063	ng/ul	76
22) Benzo(k)fluoranthene	22.96	252	2341	0.061	ng/ul#	73
23) Benzo(a)pyrene	23.50	252	2086	0.060	ng/ul#	67
24) Indeno(1,2,3-cd)pyrene	25.89	276	2838	0.062	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.90	278	2268	0.061	ng/ul#	84
26) Benzo(g,h,i)perylene	26.59	276	2620	0.063	ng/ul	91

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

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