

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN021419\
 Data File : BN004416.D
 Acq On : 14 Feb 2019 12:43
 Operator : JU/SJ
 Sample : MDL-W-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-W-01

Quant Time: Feb 14 15:00:03 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	1018	0.40	ng/ul	0.00
2) Naphthalene-d8	10.53	136	4108	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	2531	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.13	188	6950	0.40	ng/ul	0.00
16) Chrysene-d12	21.33	240	8806	0.40	ng/ul	0.00
20) Perylene-d12	23.60	264	8982	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	2316	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.16	212	7420	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.58	128	797	0.068	ng/ul#	83
5) 2-Methylnaphthalene	12.20	142	579	0.069	ng/ul	99
7) Acenaphthylene	14.11	152	721	0.062	ng/ul#	85
8) Acenaphthene	14.45	153	667	0.066	ng/ul	96
9) Fluorene	15.43	166	816	0.068	ng/ul#	97
11) Pentachlorophenol	16.78	266	37	0.025	ng/ul	94
12) Phenanthrene	17.17	178	1535	0.068	ng/ul#	92
13) Anthracene	17.27	178	1208	0.063	ng/ul#	88
15) Fluoranthene	19.20	202	1894	0.068	ng/ul	95
17) Pyrene	19.56	202	1914	0.068	ng/ul#	90
18) Benzo(a)anthracene	21.31	228	1884	0.066	ng/ul	96
19) Chrysene	21.37	228	2335	0.071	ng/ul	97
21) Benzo(b)fluoranthene	22.92	252	2513	0.070	ng/ul	75
22) Benzo(k)fluoranthene	22.96	252	2321	0.066	ng/ul#	74
23) Benzo(a)pyrene	23.50	252	2079	0.065	ng/ul#	66
24) Indeno(1,2,3-cd)pyrene	25.89	276	2870	0.068	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.90	278	2292	0.067	ng/ul#	81
26) Benzo(g,h,i)perylene	26.60	276	2674	0.070	ng/ul	91

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

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