

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN021119\
 Data File : BN004378.D
 Acq On : 11 Feb 2019 15:12
 Operator : JU/SJ
 Sample : K1410-03DL 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2F73DL

Quant Time: Feb 11 16:31:21 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	1117	0.40	ng/ul	0.00
2) Naphthalene-d8	10.54	136	4247	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.39	164	2453	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.14	188	6774	0.40	ng/ul	0.00
16) Chrysene-d12	21.33	240	7757	0.40	ng/ul	0.00
20) Perylene-d12	23.60	264	8956	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.13	152	1258	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.17	212	3896	0.18	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.59	128	4134	0.339	ng/ul	98
5) 2-Methylnaphthalene	12.20	142	1987	0.228	ng/ul	99
9) Fluorene	15.44	166	3685	0.317	ng/ul	98
11) Pentachlorophenol	16.78	266	2192	1.520	ng/ul	96
13) Anthracene	17.27	178	7959	0.429	ng/ul	100
15) Fluoranthene	19.20	202	9122	0.338	ng/ul	99
17) Pyrene	19.56	202	10517	0.421	ng/ul	99
19) Chrysene	21.37	228	8479	0.292	ng/ul	99
21) Benzo(b)fluoranthene	22.91	252	9334	0.261	ng/ul	95
23) Benzo(a)pyrene	23.50	252	11355	0.355	ng/ul	98
25) Dibenzo(a,h)anthracene	25.91	278	24546	0.720	ng/ul	98

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

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