

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061119\
 Data File : BN006261.D
 Acq On : 12 Jun 2019 00:25
 Operator : JU/SJ
 Sample : K3215-05
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DB8H9

Quant Time: Jun 12 01:12:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 05:10:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	5433	0.40	ng/ul	0.01
2) Naphthalene-d8	10.47	136	30500	0.40	ng/ul	0.04
6) Acenaphthene-d10	14.35	164	30903	0.40	ng/ul	0.06
10) Phenanthrene-d10	17.20	188	2167836	0.40	ng/ul	0.15
16) Chrysene-d12	21.43	240	310	0.40	ng/ul	0.17
20) Perylene-d12	23.49	264	22580	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	12566	0.27	ng/ul	0.04
14) Fluoranthene-d10	19.26	212	1973	0.00	ng/ul	0.17
Target Compounds				Ovalue		
3) Naphthalene	10.53	128	1915656	22.761	ng/ul#	95
5) 2-Methylnaphthalene	12.14	142	1137830	19.399	ng/ul	89
7) Acenaphthylene	14.06	152	30353	0.177	ng/ul#	1
8) Acenaphthene	14.44	153	62077	0.521	ng/ul#	51
9) Fluorene	15.39	166	58310	0.418	ng/ul#	73
17) Pyrene	19.65	202	12178	7.458	ng/ul#	1
18) Benzo(a)anthracene	21.46	228	258	0.188	ng/ul#	1
19) Chrysene	21.46	228	827	0.643	ng/ul#	1
23) Benzo(a)pyrene	23.36	252	6421	0.073	ng/ul#	2

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

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