

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN073020\
 Data File : BN011295.D
 Acq On : 30 Jul 2020 12:29
 Operator : CG/JU
 Sample : SSTDO.106
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDO.106

Quant Time: Jul 30 14:19:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 30 14:13:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	1214	0.40	ng/ul	0.00
2) Naphthalene-d8	10.22	136	4150	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	2571	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	5878	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	5384	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	4866	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.82	152	1027	0.10	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	2921	0.12	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.27	128	1246	0.100	ng/ul#	85
5) 2-Methylnaphthalene	11.90	142	811	0.100	ng/ul	98
7) Acenaphthylene	13.82	152	1210	0.097	ng/ul#	91
8) Acenaphthene	14.17	153	972	0.099	ng/ul	96
9) Fluorene	15.17	166	1142	0.103	ng/ul#	98
12) Phenanthrene	16.90	178	1966	0.101	ng/ul#	91
13) Anthracene	17.00	178	1629	0.102	ng/ul#	90
15) Fluoranthene	18.93	202	2464	0.114	ng/ul	91
17) Pyrene	19.30	202	2579	0.090	ng/ul#	93
18) Benzo(a)anthracene	21.07	228	2043	0.107	ng/ul	95
19) Chrysene	21.12	228	2438	0.096	ng/ul	98
21) Benzo(b)fluoranthene	22.58	252	1985	0.094	ng/ul#	62
22) Benzo(k)fluoranthene	22.62	252	2264	0.094	ng/ul#	56
23) Benzo(a)pyrene	23.12	252	1720	0.097	ng/ul#	39
24) Indeno(1,2,3-cd)pyrene	25.29	276	1737	0.083	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.31	278	1375	0.088	ng/ul#	53
26) Benzo(a,h,i)perylene	25.91	276	1627	0.079	ng/ul#	79

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

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