

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060320\
 Data File : BN010750.D
 Acq On : 03 Jun 2020 13:45
 Operator : CG/JU
 Sample : SSTD0.201
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.201

Quant Time: Jun 03 14:58:14 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN060320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 03 14:56:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1668	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5459	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	2776	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	5425	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	4385	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	4129	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	2438	0.27	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	4444	0.28	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	3098	0.184	ng/ul	97
5) 2-Methylnaphthalene	12.00	142	1933	0.179	ng/ul	100
7) Acenaphthylene	13.91	152	2395	0.180	ng/ul	95
8) Acenaphthene	14.26	153	2013	0.184	ng/ul	96
9) Fluorene	15.25	166	2203	0.183	ng/ul	98
11) Pentachlorophenol	16.61	266	234	0.162	ng/ul	96
12) Phenanthrene	16.98	178	3401	0.182	ng/ul	97
13) Anthracene	17.08	178	2757	0.178	ng/ul	94
15) Fluoranthene	19.01	202	3698	0.180	ng/ul	99
17) Pyrene	19.37	202	3761	0.197	ng/ul	97
18) Benzo(a)anthracene	21.13	228	2958	0.178	ng/ul	98
19) Chrysene	21.18	228	3980	0.198	ng/ul	97
21) Benzo(b)fluoranthene	22.66	252	3440	0.184	ng/ul	86
22) Benzo(k)fluoranthene	22.71	252	3742	0.187	ng/ul#	84
23) Benzo(a)pyrene	23.21	252	3010	0.187	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.43	276	4076	0.184	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.44	278	3265	0.181	ng/ul#	86
26) Benzo(g,h,i)perylene	26.07	276	3804	0.195	ng/ul	96

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

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