

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN081220\
 Data File : BN011414.D
 Acq On : 12 Aug 2020 10:50
 Operator : CG/JU
 Sample : SSTD0.203
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.203

Quant Time: Aug 12 11:59:25 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN081220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 12 11:51:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	958	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	2798	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.08	164	1462	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	3145	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	3326	0.40	ng/ul	0.00
20) Perylene-d12	23.17	264	2995	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	733	0.13	ng/ul	0.01
14) Fluoranthene-d10	18.88	212	1854	0.13	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.25	128	1654	0.222	ng/ul#	91
5) 2-Methylnaphthalene	11.88	142	963	0.219	ng/ul	100
7) Acenaphthylene	13.80	152	1352	0.204	ng/ul#	93
8) Acenaphthene	14.14	153	1085	0.206	ng/ul	98
9) Fluorene	15.15	166	1234	0.201	ng/ul#	96
11) Pentachlorophenol	16.52	266	125	0.179	ng/ul	92
12) Phenanthrene	16.88	178	2026	0.210	ng/ul	95
13) Anthracene	16.98	178	1646	0.202	ng/ul	93
15) Fluoranthene	18.91	202	2579	0.216	ng/ul	95
17) Pyrene	19.27	202	2351	0.180	ng/ul#	96
18) Benzo(a)anthracene	21.04	228	2017	0.194	ng/ul	97
19) Chrysene	21.09	228	3041	0.227	ng/ul	97
21) Benzo(b)fluoranthene	22.55	252	2372	0.210	ng/ul	87
22) Benzo(k)fluoranthene	22.59	252	2632	0.210	ng/ul#	83
23) Benzo(a)pyrene	23.08	252	2091	0.206	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.22	276	2618	0.194	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.24	278	2114	0.192	ng/ul#	81
26) Benzo(g,h,i)perylene	25.84	276	2452	0.188	ng/ul#	90

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

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