

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112020\
 Data File : BN012656.D
 Acq On : 20 Nov 2020 12:26
 Operator : CG/JU
 Sample : SST00.206
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.206

Quant Time: Nov 20 13:45:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 20 13:42:54 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	726	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	2649	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	1436	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	2937	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	2660	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	2466	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.82	152	712	0.18	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	1449	0.18	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.26	128	1588	0.208	ng/ul	93
5) 2-Methylnaphthalene	11.90	142	983	0.202	ng/ul	98
7) Acenaphthylene	13.82	152	1357	0.218	ng/ul	95
8) Acenaphthene	14.17	153	1176	0.223	ng/ul	98
9) Fluorene	15.17	166	1256	0.214	ng/ul#	96
11) Pentachlorophenol	16.52	266	182	0.214	ng/ul	92
12) Phenanthrene	16.90	178	1979	0.211	ng/ul	96
13) Anthracene	16.99	178	1702	0.212	ng/ul	95
15) Fluoranthene	18.93	202	2240	0.205	ng/ul	95
17) Pyrene	19.29	202	2384	0.199	ng/ul#	95
18) Benzo(a)anthracene	21.06	228	1809	0.184	ng/ul	98
19) Chrysene	21.11	228	2471	0.206	ng/ul	97
21) Benzo(b)fluoranthene	22.57	252	2146	0.215	ng/ul	84
22) Benzo(k)fluoranthene	22.61	252	2577	0.234	ng/ul#	85
23) Benzo(a)pyrene	23.10	252	2050	0.212	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.28	276	2623	0.211	ng/ul	98
25) Dibenzo(a,h)anthracene	25.29	278	2110	0.212	ng/ul#	86
26) Benzo(g,h,i)perylene	25.91	276	2434	0.213	ng/ul	93

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

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