

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN080620\
 Data File : BN011390.D
 Acq On : 06 Aug 2020 21:37
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.420

Quant Time: Aug 07 02:06:44 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 : Fri Aug 07 02:04:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	1224	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	4269	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	2180	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	4658	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	4993	0.40	ng/ul	0.00
20) Perylene-d12	23.17	264	4610	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.80	152	3558	0.33	ng/ul	0.00
14) Fluoranthene-d10	18.88	212	8717	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.25	128	4660	0.367	ng/ul	99
5) 2-Methylnaphthalene	11.88	142	2936	0.345	ng/ul	99
7) Acenaphthylene	13.80	152	4202	0.404	ng/ul	99
8) Acenaphthene	14.15	153	3033	0.367	ng/ul	99
9) Fluorene	15.15	166	3473	0.357	ng/ul	99
11) Pentachlorophenol	16.51	266	457	0.400	ng/ul	98
12) Phenanthrene	16.89	178	5582	0.362	ng/ul	98
13) Anthracene	16.98	178	4891	0.363	ng/ul	98
15) Fluoranthene	18.91	202	7315	0.384	ng/ul	98
17) Pyrene	19.28	202	7285	0.321	ng/ul	98
18) Benzo(a)anthracene	21.04	228	6491	0.338	ng/ul	99
19) Chrysene	21.10	228	7495	0.341	ng/ul	99
21) Benzo(b)fluoranthene	22.55	252	6732	0.350	ng/ul	96
22) Benzo(k)fluoranthene	22.59	252	7481	0.347	ng/ul	97
23) Benzo(a)pyrene	23.09	252	6158	0.368	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.24	276	7976	0.454	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.25	278	6476	0.462	ng/ul	99
26) Benzo(g,h,i)perylene	25.86	276	6977	0.439	ng/ul	99

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

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