

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120320\
 Data File : BN012845.D
 Acq On : 03 Dec 2020 13:38
 Operator : CG/JU
 Sample : SST00.803
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.803

Quant Time: Dec 03 14:10:43 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 03 13:32:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.41	152	468	0.40	ng/ul	0.00
2) Naphthalene-d8	10.17	136	1807	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.07	164	1040	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.82	188	2261	0.40	ng/ul	0.00
16) Chrysene-d12	21.04	240	1989	0.40	ng/ul	0.00
20) Perylene-d12	23.15	264	1899	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.77	152	2219	0.85	ng/ul	0.00
14) Fluoranthene-d10	18.87	212	5049	0.85	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.22	128	4203	0.773	ng/ul	93
5) 2-Methylnaphthalene	11.85	142	2797	0.805	ng/ul	99
7) Acenaphthylene	13.78	152	4057	0.800	ng/ul	96
8) Acenaphthene	14.13	153	3316	0.776	ng/ul	100
9) Fluorene	15.12	166	3855	0.820	ng/ul	99
11) Pentachlorophenol	16.48	266	526	0.733	ng/ul	95
12) Phenanthrene	16.86	178	6232	0.792	ng/ul	98
13) Anthracene	16.96	178	5538	0.810	ng/ul	96
15) Fluoranthene	18.89	202	7367	0.811	ng/ul	96
17) Pyrene	19.26	202	7366	0.828	ng/ul	98
18) Benzo(a)anthracene	21.03	228	6738	0.945	ng/ul	98
19) Chrysene	21.08	228	7536	0.846	ng/ul	99
21) Benzo(b)fluoranthene	22.52	252	7374	0.844	ng/ul	92
22) Benzo(k)fluoranthene	22.57	252	8090	0.803	ng/ul#	92
23) Benzo(a)pyrene	23.06	252	6655	0.809	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.21	276	8949	0.832	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.22	278	7158	0.829	ng/ul	92
26) Benzo(g,h,i)perylene	25.84	276	8030	0.830	ng/ul	95

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

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