

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031620\
 Data File : BN010226.D
 Acq On : 16 Mar 2020 13:51
 Operator : CG/JU
 Sample : SST00.801
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.801

Quant Time: Mar 16 14:44:44 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN031620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 16 13:58:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1801	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	5644	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3255	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	7566	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	7093	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	6724	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	7690	0.84	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	19229	0.83	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.44	128	14381	0.847	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	9707	0.833	ng/ul	100
7) Acenaphthylene	13.97	152	14138	0.851	ng/ul	99
8) Acenaphthene	14.32	153	10819	0.810	ng/ul	99
9) Fluorene	15.31	166	12937	0.843	ng/ul	99
11) Pentachlorophenol	16.65	266	1203	0.581	ng/ul	98
12) Phenanthrene	17.04	178	22146	0.847	ng/ul	99
13) Anthracene	17.13	178	19760	0.853	ng/ul	99
15) Fluoranthene	19.06	202	27280	0.881	ng/ul	98
17) Pyrene	19.42	202	26868	0.789	ng/ul	99
18) Benzo(a)anthracene	21.18	228	24439	0.786	ng/ul	99
19) Chrysene	21.23	228	27507	0.818	ng/ul	100
21) Benzo(b)fluoranthene	22.72	252	25050	0.840	ng/ul	97
22) Benzo(k)fluoranthene	22.76	252	27012	0.795	ng/ul	97
23) Benzo(a)pyrene	23.27	252	21610	0.724	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.52	276	27783	0.834	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.54	278	22847	0.869	ng/ul	98
26) Benzo(g,h,i)perylene	26.17	276	24305	0.833	ng/ul	99

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

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