

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031620\
 Data File : BN010227.D
 Acq On : 16 Mar 2020 14:26
 Operator : CG/JU
 Sample : SSTD1.602
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.602

Quant Time: Mar 16 15:18:39 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	1816	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	5725	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3398	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	7835	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	7591	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	6988	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	16037	1.72	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	41680	1.74	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.44	128	29799	1.731	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	20357	1.721	ng/ul	99
7) Acenaphthylene	13.97	152	30694	1.770	ng/ul	98
8) Acenaphthene	14.32	153	22973	1.648	ng/ul	100
9) Fluorene	15.31	166	27906	1.741	ng/ul	97
11) Pentachlorophenol	16.65	266	2809	1.310	ng/ul	98
12) Phenanthrene	17.04	178	47266	1.746	ng/ul	99
13) Anthracene	17.13	178	43265	1.803	ng/ul	99
15) Fluoranthene	19.06	202	59111	1.844	ng/ul	99
17) Pyrene	19.42	202	58030	1.592	ng/ul	98
18) Benzo(a)anthracene	21.18	228	54566	1.640	ng/ul	99
19) Chrysene	21.23	228	58097	1.613	ng/ul	100
21) Benzo(b)fluoranthene	22.72	252	53469	1.726	ng/ul	97
22) Benzo(k)fluoranthene	22.77	252	57204	1.620	ng/ul	97
23) Benzo(a)pyrene	23.27	252	46669	1.504	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.53	276	59675	1.724	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.54	278	49137	1.797	ng/ul	97
26) Benzo(g,h,i)perylene	26.17	276	51313	1.692	ng/ul	99

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

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