

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN081220\
 Data File : BN011418.D
 Acq On : 12 Aug 2020 13:13
 Operator : CG/JU
 Sample : SSTD3.201
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.201

Quant Time: Aug 12 13:55:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN081220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 12 11:51:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	788	0.40	ng/ul	0.00
2) Naphthalene-d8	10.19	136	2225	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.07	164	1179	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.83	188	2876	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	3377	0.40	ng/ul	0.00
20) Perylene-d12	23.17	264	2339	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.79	152	11402	2.54	ng/ul	-0.01
14) Fluoranthene-d10	18.87	212	24698	1.96	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.24	128	18880	3.188	ng/ul	97
5) 2-Methylnaphthalene	11.86	142	15271	4.369	ng/ul	99
7) Acenaphthylene	13.79	152	18708	3.502	ng/ul	97
8) Acenaphthene	14.14	153	14535	3.423	ng/ul	98
9) Fluorene	15.13	166	17004	3.433	ng/ul	97
11) Pentachlorophenol	16.49	266	2251	3.528	ng/ul	98
12) Phenanthrene	16.87	178	33436	3.787	ng/ul	96
13) Anthracene	16.97	178	28728	3.860	ng/ul	97
15) Fluoranthene	18.90	202	34165	3.127	ng/ul	96
17) Pyrene	19.27	202	33514	2.525	ng/ul	96
18) Benzo(a)anthracene	21.03	228	40133	3.804	ng/ul	98
19) Chrysene	21.09	228	44186	3.244	ng/ul	99
21) Benzo(b)fluoranthene	22.54	252	37917	4.293	ng/ul	91
22) Benzo(k)fluoranthene	22.59	252	40614	4.146	ng/ul#	91
23) Benzo(a)pyrene	23.07	252	29093	3.678	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.22	276	38847	3.677	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.23	278	32298	3.757	ng/ul	93
26) Benzo(g,h,i)perylene	25.84	276	34320	3.370	ng/ul	95

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

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