

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN072020\
 Data File : BN011206.D
 Acq On : 20 Jul 2020 12:21
 Operator : CG/JU
 Sample : SSTD3.256
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.204

Quant Time: Jul 20 12:51:32 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN072020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 20 11:25:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	1260	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	4585	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	2686	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.87	188	5762	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	4758	0.40	ng/ul	0.00
20) Perylene-d12	23.23	264	3611	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	34152	3.17	ng/ul	0.00
14) Fluoranthene-d10	18.91	212	82409	3.31	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.28	128	39381	2.910	ng/ul	96
5) 2-Methylnaphthalene	11.91	142	27134	3.199	ng/ul	100
7) Acenaphthylene	13.83	152	39946	3.336	ng/ul	97
8) Acenaphthene	14.18	153	30194	3.009	ng/ul	99
9) Fluorene	15.18	166	35595	3.227	ng/ul	99
11) Pentachlorophenol	16.54	266	4388	3.252	ng/ul	96
12) Phenanthrene	16.92	178	57597	3.072	ng/ul	98
13) Anthracene	17.00	178	53012	3.423	ng/ul	96
15) Fluoranthene	18.94	202	70990	3.374	ng/ul	98
17) Pyrene	19.31	202	68066	3.292	ng/ul	99
18) Benzo(a)anthracene	21.08	228	54913	3.251	ng/ul	96
19) Chrysene	21.13	228	59011	2.747	ng/ul	98
21) Benzo(b)fluoranthene	22.59	252	49848	3.099	ng/ul	85
22) Benzo(k)fluoranthene	22.64	252	52460	3.081	ng/ul#	87
23) Benzo(a)pyrene	23.13	252	41062	3.023	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.31	276	49593	2.665	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.32	278	39469	2.597	ng/ul#	85
26) Benzo(g,h,i)perylene	25.93	276	44508	2.585	ng/ul	96

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

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