

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN012987.D
 Acq On : 08 Dec 2020 22:41
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.432

Quant Time: Dec 09 00:15:02 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 : Wed Dec 09 00:11:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1046	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	4270	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	2359	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	4701	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3151	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	2852	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	2549	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	4729	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	4977	0.403	ng/ul	93
5) 2-Methylnaphthalene	11.83	142	3258	0.401	ng/ul	99
7) Acenaphthylene	13.76	152	4683	0.419	ng/ul	97
8) Acenaphthene	14.11	153	3689	0.403	ng/ul	97
9) Fluorene	15.11	166	4063	0.387	ng/ul	97
11) Pentachlorophenol	16.47	266	507	0.372	ng/ul	97
12) Phenanthrene	16.85	178	6330	0.402	ng/ul	98
13) Anthracene	16.94	178	5685	0.424	ng/ul	97
15) Fluoranthene	18.88	202	6861	0.366	ng/ul	99
17) Pyrene	19.25	202	6980	0.469	ng/ul	99
18) Benzo(a)anthracene	21.01	228	5274	0.405	ng/ul	98
19) Chrysene	21.06	228	5930	0.412	ng/ul	99
21) Benzo(b)fluoranthene	22.51	252	5444	0.413	ng/ul	96
22) Benzo(k)fluoranthene	22.55	252	5984	0.410	ng/ul	97
23) Benzo(a)pyrene	23.04	252	5140	0.420	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.18	276	6071	0.378	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.20	278	4896	0.385	ng/ul	94
26) Benzo(g,h,i)perylene	25.81	276	5246	0.361	ng/ul	98

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

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