

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011720\
 Data File : BN009396.D
 Acq On : 17 Jan 2020 15:20
 Operator : JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.421

Quant Time: Jan 17 15:50:59 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN011020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 17 12:28:30 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	1223	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	4068	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	2273	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	4643	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	3667	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	3595	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	2585	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	5459	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.27	128	4965	0.406	ng/ul	100
5) 2-Methylnaphthalene	11.91	142	3250	0.387	ng/ul	100
7) Acenaphthylene	13.82	152	4239	0.365	ng/ul	98
8) Acenaphthene	14.17	153	3550	0.381	ng/ul	100
9) Fluorene	15.17	166	3915	0.365	ng/ul	98
11) Pentachlorophenol	16.55	266	375	0.295	ng/ul	97
12) Phenanthrene	16.90	178	6389	0.398	ng/ul	100
13) Anthracene	17.00	178	5396	0.379	ng/ul	99
15) Fluoranthene	18.93	202	7226	0.380	ng/ul	98
17) Pyrene	19.29	202	7407	0.421	ng/ul	99
18) Benzo(a)anthracene	21.05	228	6011	0.374	ng/ul	99
19) Chrysene	21.10	228	7381	0.424	ng/ul	99
21) Benzo(b)fluoranthene	22.56	252	6407	0.402	ng/ul	96
22) Benzo(k)fluoranthene	22.60	252	7635	0.420	ng/ul	97
23) Benzo(a)pyrene	23.10	252	6136	0.384	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.26	276	6776	0.381	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.27	278	5322	0.378	ng/ul	94
26) Benzo(g,h,i)perylene	25.88	276	6169	0.395	ng/ul	98

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

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