

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090120\
 Data File : BN011681.D
 Acq On : 01 Sep 2020 13:00
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.408

Quant Time: Sep 01 13:37:59 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 : Tue Sep 01 13:14:23 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	1303	0.40	ng/ul	0.00
2) Naphthalene-d8	10.53	136	4656	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	2531	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.13	188	5694	0.40	ng/ul	0.00
16) Chrysene-d12	21.33	240	4890	0.40	ng/ul	0.00
20) Perylene-d12	23.58	264	4457	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	3014	0.46	ng/ul	0.00
14) Fluoranthene-d10	19.16	212	6408	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.58	128	5529	0.407	ng/ul	98
5) 2-Methylnaphthalene	12.20	142	3724	0.434	ng/ul	98
7) Acenaphthylene	14.10	152	4954	0.389	ng/ul	99
8) Acenaphthene	14.45	153	3995	0.403	ng/ul	98
9) Fluorene	15.43	166	4641	0.397	ng/ul	99
11) Pentachlorophenol	16.78	266	685	0.507	ng/ul	98
12) Phenanthrene	17.17	178	7728	0.401	ng/ul	97
13) Anthracene	17.26	178	6978	0.440	ng/ul	99
15) Fluoranthene	19.19	202	8856	0.400	ng/ul	99
17) Pyrene	19.55	202	8879	0.465	ng/ul	98
18) Benzo(a)anthracene	21.31	228	7685	0.468	ng/ul	98
19) Chrysene	21.36	228	8671	0.392	ng/ul	99
21) Benzo(b)fluoranthene	22.90	252	8580	0.443	ng/ul	97
22) Benzo(k)fluoranthene	22.95	252	8485	0.396	ng/ul	98
23) Benzo(a)pyrene	23.48	252	7148	0.439	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.85	276	9461	0.430	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.87	278	7834	0.437	ng/ul	98
26) Benzo(g,h,i)perylene	26.54	276	8242	0.400	ng/ul	99

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

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