

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090120\
 Data File : BN011687.D
 Acq On : 01 Sep 2020 16:39
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.409

Quant Time: Sep 01 17:09:05 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	7.74	152	1674	0.40	ng/ul	0.00
2) Naphthalene-d8	10.53	136	6060	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	3190	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.13	188	6795	0.40	ng/ul	0.00
16) Chrysene-d12	21.32	240	5784	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	5273	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	4057	0.47	ng/ul	0.00
14) Fluoranthene-d10	19.16	212	7856	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.58	128	7604	0.430	ng/ul	98
5) 2-Methylnaphthalene	12.19	142	5099	0.457	ng/ul	97
7) Acenaphthylene	14.10	152	6505	0.405	ng/ul	98
8) Acenaphthene	14.44	153	5227	0.418	ng/ul	98
9) Fluorene	15.43	166	6008	0.408	ng/ul	98
11) Pentachlorophenol	16.78	266	853	0.529	ng/ul	99
12) Phenanthrene	17.17	178	9777	0.425	ng/ul	97
13) Anthracene	17.26	178	8830	0.467	ng/ul	97
15) Fluoranthene	19.19	202	11071	0.419	ng/ul	99
17) Pyrene	19.55	202	10996	0.486	ng/ul	98
18) Benzo(a)anthracene	21.31	228	9592	0.494	ng/ul	98
19) Chrysene	21.36	228	10472	0.400	ng/ul	99
21) Benzo(b)fluoranthene	22.90	252	10318	0.450	ng/ul	97
22) Benzo(k)fluoranthene	22.94	252	10960	0.433	ng/ul	97
23) Benzo(a)pyrene	23.47	252	8976	0.466	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.84	276	12040	0.462	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.86	278	9841	0.464	ng/ul	99
26) Benzo(g,h,i)perylene	26.53	276	10099	0.414	ng/ul	99

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

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