

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120719\
 Data File : BN008887.D
 Acq On : 07 Dec 2019 13:15
 Operator : JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.405

Quant Time: Dec 09 00:13:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 06 12:03:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2219	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	9380	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	6310	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	14076	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	12780	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	13286	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	6325	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	16635	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	11059	0.391	ng/ul	100
5) 2-Methylnaphthalene	11.99	142	8224	0.416	ng/ul	100
7) Acenaphthylene	13.90	152	12133	0.348	ng/ul	99
8) Acenaphthene	14.25	153	9472	0.366	ng/ul	100
9) Fluorene	15.24	166	11247	0.370	ng/ul	100
11) Pentachlorophenol	16.59	266	1498	0.281	ng/ul	98
12) Phenanthrene	16.97	178	19381	0.408	ng/ul	99
13) Anthracene	17.07	178	17488	0.390	ng/ul	99
15) Fluoranthene	19.00	202	22577	0.387	ng/ul	98
17) Pyrene	19.36	202	22713	0.443	ng/ul	99
18) Benzo(a)anthracene	21.12	228	22220	0.432	ng/ul	100
19) Chrysene	21.17	228	22236	0.440	ng/ul	99
21) Benzo(b)fluoranthene	22.65	252	22311	0.407	ng/ul	100
22) Benzo(k)fluoranthene	22.69	252	22818	0.420	ng/ul	99
23) Benzo(a)pyrene	23.19	252	20872	0.406	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.39	276	22871	0.380	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.41	278	18537	0.382	ng/ul	99
26) Benzo(g,h,i)perylene	26.03	276	18698	0.372	ng/ul	99

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

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