

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090420\
 Data File : BN011727.D
 Acq On : 05 Sep 2020 07:24
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.412

Quant Time: Sep 05 07:54:37 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 : Sat Sep 05 05:40:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2772	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	10791	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	6036	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	12956	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	10933	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	10204	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	7107	0.47	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	14407	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	12672	0.403	ng/ul	97
5) 2-Methylnaphthalene	12.19	142	8473	0.426	ng/ul	100
7) Acenaphthylene	14.09	152	11558	0.380	ng/ul	97
8) Acenaphthene	14.44	153	9248	0.391	ng/ul	98
9) Fluorene	15.43	166	10583	0.380	ng/ul	99
11) Pentachlorophenol	16.77	266	1476	0.480	ng/ul	98
12) Phenanthrene	17.16	178	17271	0.394	ng/ul	97
13) Anthracene	17.25	178	15621	0.433	ng/ul	96
15) Fluoranthene	19.18	202	19476	0.386	ng/ul	96
17) Pyrene	19.54	202	19594	0.459	ng/ul#	94
18) Benzo(a)anthracene	21.30	228	17430	0.475	ng/ul	97
19) Chrysene	21.35	228	18516	0.374	ng/ul	99
21) Benzo(b)fluoranthene	22.88	252	18562	0.419	ng/ul	96
22) Benzo(k)fluoranthene	22.93	252	18855	0.385	ng/ul	96
23) Benzo(a)pyrene	23.46	252	16703	0.448	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.82	276	20971	0.416	ng/ul	94
25) Dibenzo(a,h)anthracene	25.84	278	16943	0.412	ng/ul	96
26) Benzo(g,h,i)perylene	26.51	276	17424	0.369	ng/ul	98

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

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