

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN012972.D
 Acq On : 08 Dec 2020 12:45
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.431

Quant Time: Dec 08 13:43:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 07 12:24:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1101	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	4429	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	2390	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	4750	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3570	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	3291	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	2640	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	4972	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	5080	0.396	ng/ul#	92
5) 2-Methylnaphthalene	11.83	142	3375	0.401	ng/ul	99
7) Acenaphthylene	13.76	152	4749	0.420	ng/ul	96
8) Acenaphthene	14.11	153	3673	0.396	ng/ul	100
9) Fluorene	15.11	166	4054	0.381	ng/ul	98
11) Pentachlorophenol	16.47	266	563	0.409	ng/ul	95
12) Phenanthrene	16.85	178	6385	0.401	ng/ul	98
13) Anthracene	16.95	178	5751	0.424	ng/ul	96
15) Fluoranthene	18.88	202	7250	0.383	ng/ul	100
17) Pyrene	19.25	202	7381	0.438	ng/ul	99
18) Benzo(a)anthracene	21.02	228	6079	0.412	ng/ul	98
19) Chrysene	21.07	228	6663	0.408	ng/ul	99
21) Benzo(b)fluoranthene	22.52	252	6151	0.404	ng/ul	96
22) Benzo(k)fluoranthene	22.56	252	6625	0.394	ng/ul	98
23) Benzo(a)pyrene	23.05	252	5639	0.399	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.19	276	6680	0.360	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.20	278	5390	0.367	ng/ul	96
26) Benzo(g,h,i)perylene	25.82	276	5663	0.337	ng/ul	98

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

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