

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112420\
 Data File : BN012729.D
 Acq On : 24 Nov 2020 22:51
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.414

Quant Time: Nov 25 01:53:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 25 01:45:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	1084	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	4449	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	2418	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	4696	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	4191	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	4363	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.80	152	2609	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	4639	0.38	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.25	128	5476	0.409	ng/ul	97
5) 2-Methylnaphthalene	11.88	142	3730	0.436	ng/ul	99
7) Acenaphthylene	13.81	152	5286	0.448	ng/ul	97
8) Acenaphthene	14.16	153	3980	0.400	ng/ul	99
9) Fluorene	15.15	166	4373	0.400	ng/ul	100
11) Pentachlorophenol	16.51	266	767	0.514	ng/ul	96
12) Phenanthrene	16.89	178	6629	0.406	ng/ul	98
13) Anthracene	16.98	178	6048	0.426	ng/ul	96
15) Fluoranthene	18.92	202	8110	0.430	ng/ul	98
17) Pyrene	19.28	202	8269	0.441	ng/ul	98
18) Benzo(a)anthracene	21.05	228	8177	0.544	ng/ul	98
19) Chrysene	21.10	228	9387	0.500	ng/ul	99
21) Benzo(b)fluoranthene	22.55	252	9709	0.484	ng/ul	95
22) Benzo(k)fluoranthene	22.59	252	10086	0.436	ng/ul	95
23) Benzo(a)pyrene	23.08	252	8884	0.470	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.25	276	11531	0.467	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.26	278	9167	0.462	ng/ul	95
26) Benzo(g,h,i)perylene	25.89	276	9789	0.440	ng/ul	96

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

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