

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112020\
 Data File : BN012690.D
 Acq On : 21 Nov 2020 17:34
 Operator : CG/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.411

Quant Time: Nov 23 01:25:36 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 : Mon Nov 23 01:22:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	1036	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	3962	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	2033	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	3863	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	3260	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	3545	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	2230	0.39	ng/ul	-0.01
14) Fluoranthene-d10	18.89	212	3859	0.38	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.26	128	4704	0.394	ng/ul	97
5) 2-Methylnaphthalene	11.89	142	3088	0.406	ng/ul	98
7) Acenaphthylene	13.82	152	4264	0.430	ng/ul	98
8) Acenaphthene	14.16	153	3240	0.388	ng/ul	99
9) Fluorene	15.16	166	3494	0.380	ng/ul	98
11) Pentachlorophenol	16.51	266	522	0.425	ng/ul	97
12) Phenanthrene	16.89	178	5323	0.396	ng/ul	97
13) Anthracene	16.99	178	4756	0.407	ng/ul	97
15) Fluoranthene	18.92	202	5995	0.386	ng/ul	97
17) Pyrene	19.29	202	6249	0.429	ng/ul	96
18) Benzo(a)anthracene	21.05	228	5407	0.463	ng/ul	98
19) Chrysene	21.10	228	6293	0.431	ng/ul	99
21) Benzo(b)fluoranthene	22.56	252	6483	0.398	ng/ul	97
22) Benzo(k)fluoranthene	22.60	252	6748	0.359	ng/ul	97
23) Benzo(a)pyrene	23.09	252	6072	0.396	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.26	276	7880	0.393	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.27	278	6197	0.384	ng/ul	96
26) Benzo(g,h,i)perylene	25.89	276	6787	0.376	ng/ul	96

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

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