

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112420\
 Data File : BN012709.D
 Acq On : 24 Nov 2020 10:30
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.412

Quant Time: Nov 24 12:00:27 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 : Tue Nov 24 11:59:48 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	738	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	2919	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	1621	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	3286	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	3023	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	3053	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	1683	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	3315	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.25	128	3554	0.404	ng/ul	97
5) 2-Methylnaphthalene	11.88	142	2452	0.437	ng/ul	99
7) Acenaphthylene	13.81	152	3458	0.437	ng/ul	98
8) Acenaphthene	14.16	153	2713	0.407	ng/ul	99
9) Fluorene	15.15	166	2999	0.409	ng/ul	99
11) Pentachlorophenol	16.51	266	467	0.448	ng/ul	96
12) Phenanthrene	16.89	178	4698	0.411	ng/ul	98
13) Anthracene	16.98	178	4162	0.419	ng/ul	98
15) Fluoranthene	18.92	202	5766	0.437	ng/ul	100
17) Pyrene	19.28	202	5921	0.438	ng/ul	97
18) Benzo(a)anthracene	21.05	228	5652	0.522	ng/ul	99
19) Chrysene	21.10	228	6924	0.512	ng/ul	100
21) Benzo(b)fluoranthene	22.56	252	6870	0.489	ng/ul	95
22) Benzo(k)fluoranthene	22.60	252	7513	0.464	ng/ul	97
23) Benzo(a)pyrene	23.09	252	6243	0.472	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.26	276	8207	0.475	ng/ul#	56
25) Dibenzo(a,h)anthracene	25.27	278	6511	0.469	ng/ul	95
26) Benzo(g,h,i)perylene	25.89	276	7173	0.461	ng/ul	98

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

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