

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022121\
 Data File : BN013712.D
 Acq On : 21 Feb 2021 06:06
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.418

Quant Time: Feb 21 06:39:41 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN022121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 20 23:16:49 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	4065	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	15916	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	9047	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	18880	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	16932	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	15927	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	10175	0.44	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	21240	0.42	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.35	128	18521	0.441	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	12310	0.431	ng/ul	99
7) Acenaphthylene	13.90	152	15759	0.448	ng/ul	100
8) Acenaphthene	14.24	153	13192	0.435	ng/ul	100
9) Fluorene	15.23	166	14848	0.422	ng/ul	99
11) Pentachlorophenol	16.59	266	1853	0.422	ng/ul	98
12) Phenanthrene	16.97	178	24266	0.428	ng/ul	100
13) Anthracene	17.06	178	21948	0.441	ng/ul	100
15) Fluoranthene	19.00	202	27577	0.426	ng/ul	100
17) Pyrene	19.36	202	28476	0.440	ng/ul	100
18) Benzo(a)anthracene	21.12	228	25673	0.441	ng/ul	99
19) Chrysene	21.17	228	26464	0.439	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	26387	0.433	ng/ul	99
22) Benzo(k)fluoranthene	22.70	252	27293	0.435	ng/ul	100
23) Benzo(a)pyrene	23.20	252	24584	0.456	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.45	276	30146	0.433	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.46	278	24516	0.432	ng/ul	99
26) Benzo(g,h,i)perylene	26.10	276	25883	0.431	ng/ul	99

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

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