

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN013036.D
 Acq On : 10 Dec 2020 06:21
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 106 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.437

Quant Time: Dec 10 06:56:24 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 : Thu Dec 10 05:16:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	1340	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.14	136	5853	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.04	164	3219	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	6505	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	4676	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	4329	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	3538	0.41	ng/ul	-0.01
14) Fluoranthene-d10	18.85	212	6829	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.19	128	6576	0.388	ng/ul	93
5) 2-Methylnaphthalene	11.82	142	4499	0.404	ng/ul	99
7) Acenaphthylene	13.76	152	6577	0.432	ng/ul	96
8) Acenaphthene	14.11	153	5025	0.402	ng/ul	99
9) Fluorene	15.11	166	5493	0.383	ng/ul	99
11) Pentachlorophenol	16.47	266	793	0.421	ng/ul	96
12) Phenanthrene	16.84	178	8712	0.399	ng/ul	98
13) Anthracene	16.94	178	8211	0.442	ng/ul	95
15) Fluoranthene	18.88	202	9956	0.384	ng/ul	99
17) Pyrene	19.25	202	10116	0.458	ng/ul	98
18) Benzo(a)anthracene	21.01	228	8238	0.426	ng/ul	98
19) Chrysene	21.07	228	8716	0.408	ng/ul	99
21) Benzo(b)fluoranthene	22.51	252	8202	0.410	ng/ul	92
22) Benzo(k)fluoranthene	22.55	252	8492	0.384	ng/ul	96
23) Benzo(a)pyrene	23.04	252	7623	0.410	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	25.18	276	8534	0.350	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.20	278	6836	0.354	ng/ul	95
26) Benzo(g,h,i)perylene	25.81	276	7221	0.327	ng/ul	97

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

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