

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN013041.D
 Acq On : 10 Dec 2020 11:28
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 111 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.438

Quant Time: Dec 10 11:58:20 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 10:29:27 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	843	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	3589	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.04	164	1992	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	4084	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3041	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	2905	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	2250	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	4544	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.19	128	4297	0.414	ng/ul	94
5) 2-Methylnaphthalene	11.82	142	2884	0.423	ng/ul	98
7) Acenaphthylene	13.76	152	4218	0.447	ng/ul	97
8) Acenaphthene	14.11	153	3249	0.420	ng/ul	99
9) Fluorene	15.11	166	3569	0.402	ng/ul	98
11) Pentachlorophenol	16.47	266	495	0.418	ng/ul	92
12) Phenanthrene	16.84	178	5748	0.420	ng/ul	98
13) Anthracene	16.94	178	5261	0.451	ng/ul	96
15) Fluoranthene	18.88	202	6659	0.409	ng/ul	98
17) Pyrene	19.25	202	6728	0.468	ng/ul	99
18) Benzo(a)anthracene	21.01	228	5603	0.445	ng/ul	99
19) Chrysene	21.07	228	6027	0.434	ng/ul	99
21) Benzo(b)fluoranthene	22.51	252	5652	0.421	ng/ul	96
22) Benzo(k)fluoranthene	22.55	252	6098	0.411	ng/ul	98
23) Benzo(a)pyrene	23.04	252	5285	0.424	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.18	276	6187	0.378	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.19	278	4995	0.386	ng/ul	95
26) Benzo(g,h,i)perylene	25.81	276	5360	0.362	ng/ul	98

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

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