

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102020\
 Data File : BN012375.D
 Acq On : 20 Oct 2020 09:55
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.427

Quant Time: Oct 20 10:52:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 20 10:51:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	1107	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	4819	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	2696	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	5139	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	4105	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	4197	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	2991	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	5319	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.41	128	5520	0.397	ng/ul	100
5) 2-Methylnaphthalene	12.04	142	3649	0.413	ng/ul	100
7) Acenaphthylene	13.95	152	4818	0.412	ng/ul	99
8) Acenaphthene	14.30	153	4056	0.410	ng/ul	99
9) Fluorene	15.29	166	4439	0.403	ng/ul	99
11) Pentachlorophenol	16.64	266	569	0.383	ng/ul	96
12) Phenanthrene	17.03	178	6722	0.410	ng/ul	99
13) Anthracene	17.12	178	5720	0.407	ng/ul	99
15) Fluoranthene	19.05	202	7226	0.377	ng/ul	98
17) Pyrene	19.41	202	7461	0.403	ng/ul	99
18) Benzo(a)anthracene	21.17	228	6080	0.400	ng/ul	98
19) Chrysene	21.22	228	7388	0.399	ng/ul	98
21) Benzo(b)fluoranthene	22.71	252	7558	0.446	ng/ul	97
22) Benzo(k)fluoranthene	22.76	252	7900	0.421	ng/ul	99
23) Benzo(a)pyrene	23.27	252	7041	0.427	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.53	276	9034	0.426	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.54	278	7303	0.431	ng/ul	99
26) Benzo(g,h,i)perylene	26.19	276	8055	0.415	ng/ul	99

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

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