

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102219\
 Data File : BN008376.D
 Acq On : 24 Oct 2019 07:03
 Operator : JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.413

Quant Time: Oct 24 12:36:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 14:07:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	1416	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	6864	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.21	164	4496	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	9139	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	7452	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	6545	0.40	ng/ul	-0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	4401	0.38	ng/ul	-0.01
14) Fluoranthene-d10	19.00	212	11357	0.36	ng/ul	-0.01

Target Compounds				Ovalue		
3) Naphthalene	10.39	128	7722	0.395	ng/ul	99
5) 2-Methylnaphthalene	12.02	142	5240	0.391	ng/ul	99
7) Acenaphthylene	13.93	152	7637	0.357	ng/ul	99
8) Acenaphthene	14.27	153	6300	0.378	ng/ul	97
9) Fluorene	15.27	166	6917	0.366	ng/ul	96
11) Pentachlorophenol	16.65	266	537	0.243	ng/ul	99
12) Phenanthrene	17.00	178	11004	0.411	ng/ul	99
13) Anthracene	17.10	178	9596	0.408	ng/ul	99
15) Fluoranthene	19.03	202	13248	0.375	ng/ul	96
17) Pyrene	19.40	202	13628	0.459	ng/ul	97
18) Benzo(a)anthracene	21.16	228	9978	0.368	ng/ul	99
19) Chrysene	21.21	228	13141	0.395	ng/ul	99
21) Benzo(b)fluoranthene	22.70	252	9529	0.447	ng/ul	94
22) Benzo(k)fluoranthene	22.74	252	11297	0.468	ng/ul	100
23) Benzo(a)pyrene	23.25	252	9555	0.426	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.49	276	10048	0.381	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.50	278	8033	0.385	ng/ul	99
26) Benzo(g,h,i)perylene	26.14	276	8894	0.367	ng/ul	97

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

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