

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102219\
 Data File : BN008345.D
 Acq On : 23 Oct 2019 09:54
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.410

Quant Time: Oct 23 11:06:09 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 01:44:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1524	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	7649	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.22	164	5162	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	11533	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	10883	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	9621	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	5077	0.39	ng/ul	-0.01
14) Fluoranthene-d10	19.00	212	15809	0.40	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.39	128	8703	0.400	ng/ul	100
5) 2-Methylnaphthalene	12.02	142	6017	0.403	ng/ul	99
7) Acenaphthylene	13.94	152	9143	0.372	ng/ul	99
8) Acenaphthene	14.28	153	7280	0.381	ng/ul	98
9) Fluorene	15.27	166	8233	0.379	ng/ul	97
11) Pentachlorophenol	16.64	266	1011	0.362	ng/ul	99
12) Phenanthrene	17.01	178	14156	0.419	ng/ul	100
13) Anthracene	17.10	178	12910	0.434	ng/ul	100
15) Fluoranthene	19.03	202	18462	0.414	ng/ul	96
17) Pyrene	19.40	202	19092	0.441	ng/ul	98
18) Benzo(a)anthracene	21.16	228	15913	0.402	ng/ul	99
19) Chrysene	21.21	228	18745	0.386	ng/ul	99
21) Benzo(b)fluoranthene	22.70	252	14582	0.466	ng/ul	94
22) Benzo(k)fluoranthene	22.74	252	15967	0.450	ng/ul	98
23) Benzo(a)pyrene	23.25	252	14140	0.429	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.50	276	14320	0.369	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.51	278	11536	0.376	ng/ul	99
26) Benzo(g,h,i)perylene	26.15	276	12409	0.349	ng/ul	98

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

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