

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102219\
 Data File : BN008333.D
 Acq On : 23 Oct 2019 02:08
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.409

Quant Time: Oct 23 11:26:34 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	1592	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	7836	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5155	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	10485	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	8854	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	7763	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	5137	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	12999	0.36	ng/ul	0.00

Target Compounds						Ovalue
3) Naphthalene	10.39	128	8898	0.399	ng/ul	99
5) 2-Methylnaphthalene	12.02	142	6092	0.398	ng/ul	100
7) Acenaphthylene	13.94	152	9049	0.369	ng/ul	99
8) Acenaphthene	14.28	153	7217	0.378	ng/ul	98
9) Fluorene	15.28	166	8003	0.369	ng/ul	100
11) Pentachlorophenol	16.65	266	717	0.283	ng/ul	99
12) Phenanthrene	17.01	178	12786	0.417	ng/ul	100
13) Anthracene	17.11	178	11464	0.424	ng/ul	100
15) Fluoranthene	19.04	202	15160	0.374	ng/ul	97
17) Pyrene	19.40	202	15632	0.443	ng/ul	97
18) Benzo(a)anthracene	21.16	228	12385	0.385	ng/ul	99
19) Chrysene	21.21	228	15291	0.387	ng/ul	99
21) Benzo(b)fluoranthene	22.70	252	11611	0.459	ng/ul	97
22) Benzo(k)fluoranthene	22.74	252	12623	0.441	ng/ul	100
23) Benzo(a)pyrene	23.25	252	11057	0.416	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.50	276	11465	0.366	ng/ul	95
25) Dibenzo(a,h)anthracene	25.51	278	9244	0.373	ng/ul	99
26) Benzo(g,h,i)perylene	26.15	276	10013	0.349	ng/ul	97

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

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