

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101419\
 Data File : BN008149.D
 Acq On : 15 Oct 2019 00:00
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.416

Quant Time: Oct 15 03:03:02 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 : Mon Oct 14 17:19:29 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1878	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	9704	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	6456	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	15769	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	15273	0.40	ng/ul	0.00
20) Perylene-d12	23.38	264	12671	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	6236	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	19084	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	10109	0.366	ng/ul	100
5) 2-Methylnaphthalene	12.06	142	7179	0.379	ng/ul	100
7) Acenaphthylene	13.97	152	11123	0.362	ng/ul	100
8) Acenaphthene	14.32	153	9750	0.408	ng/ul	99
9) Fluorene	15.31	166	9987	0.368	ng/ul	98
11) Pentachlorophenol	16.67	266	1212	0.318	ng/ul	100
12) Phenanthrene	17.05	178	16510	0.358	ng/ul	99
13) Anthracene	17.14	178	14541	0.358	ng/ul	100
15) Fluoranthene	19.06	202	20997	0.344	ng/ul	98
17) Pyrene	19.43	202	24194	0.398	ng/ul	99
18) Benzo(a)anthracene	21.18	228	19649	0.354	ng/ul	99
19) Chrysene	21.24	228	23709	0.348	ng/ul	100
21) Benzo(b)fluoranthene	22.73	252	18273	0.443	ng/ul	97
22) Benzo(k)fluoranthene	22.78	252	20379	0.436	ng/ul	100
23) Benzo(a)pyrene	23.29	252	17650	0.407	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.56	276	17952	0.351	ng/ul	98
25) Dibenzo(a,h)anthracene	25.57	278	14295	0.354	ng/ul	100
26) Benzo(g,h,i)perylene	26.21	276	15848	0.338	ng/ul	98

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

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