

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112019\
 Data File : BN008653.D
 Acq On : 20 Nov 2019 10:08
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.415

Quant Time: Nov 20 11:39:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 20 04:45:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	3622	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	13053	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	7273	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	14524	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	12307	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	12470	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	8266	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	16929	0.38	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.43	128	15196	0.386	ng/ul	99
5) 2-Methylnaphthalene	12.04	142	10477	0.381	ng/ul	100
7) Acenaphthylene	13.96	152	14556	0.362	ng/ul	99
8) Acenaphthene	14.31	153	10717	0.359	ng/ul	99
9) Fluorene	15.30	166	12149	0.347	ng/ul	99
11) Pentachlorophenol	16.65	266	1793	0.326	ng/ul	98
12) Phenanthrene	17.03	178	19737	0.403	ng/ul	99
13) Anthracene	17.11	178	18331	0.396	ng/ul	99
15) Fluoranthene	19.05	202	22485	0.374	ng/ul	94
17) Pyrene	19.41	202	22737	0.461	ng/ul	99
18) Benzo(a)anthracene	21.16	228	21674	0.438	ng/ul	99
19) Chrysene	21.22	228	21150	0.435	ng/ul	99
21) Benzo(b)fluoranthene	22.71	252	20863	0.406	ng/ul	94
22) Benzo(k)fluoranthene	22.75	252	20369	0.399	ng/ul	98
23) Benzo(a)pyrene	23.25	252	19673	0.408	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.49	276	22283	0.395	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.50	278	17895	0.393	ng/ul	99
26) Benzo(g,h,i)perylene	26.14	276	19053	0.403	ng/ul	97

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

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