

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111919\
 Data File : BN008624.D
 Acq On : 19 Nov 2019 11:25
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.411

Quant Time: Nov 19 12:36:25 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 : Tue Nov 19 01:47:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2991	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	11139	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	6324	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	12868	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	10882	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	11281	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	7414	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	15997	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.43	128	13361	0.398	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	9446	0.403	ng/ul	100
7) Acenaphthylene	13.97	152	13920	0.398	ng/ul	98
8) Acenaphthene	14.31	153	9818	0.378	ng/ul	99
9) Fluorene	15.30	166	11115	0.365	ng/ul	99
11) Pentachlorophenol	16.65	266	1744	0.358	ng/ul	98
12) Phenanthrene	17.03	178	17979	0.414	ng/ul	99
13) Anthracene	17.13	178	17421	0.425	ng/ul	99
15) Fluoranthene	19.06	202	20928	0.393	ng/ul	98
17) Pyrene	19.42	202	21568	0.494	ng/ul	100
18) Benzo(a)anthracene	21.17	228	20006	0.457	ng/ul	98
19) Chrysene	21.23	228	18926	0.440	ng/ul	99
21) Benzo(b)fluoranthene	22.72	252	18962	0.408	ng/ul	86
22) Benzo(k)fluoranthene	22.76	252	18775	0.407	ng/ul#	80
23) Benzo(a)pyrene	23.27	252	18237	0.418	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.52	276	21264	0.416	ng/ul	99
25) Dibenzo(a,h)anthracene	25.53	278	17092	0.414	ng/ul#	91
26) Benzo(g,h,i)perylene	26.16	276	17806	0.417	ng/ul	93

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

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