

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111919\
 Data File : BN008632.D
 Acq On : 19 Nov 2019 16:14
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.412

Quant Time: Nov 19 16:53:28 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	3255	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	12078	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	6884	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	14103	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	11962	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	12144	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	8133	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	17658	0.41	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.43	128	14549	0.400	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	10266	0.404	ng/ul	100
7) Acenaphthylene	13.96	152	15003	0.394	ng/ul	99
8) Acenaphthene	14.31	153	10678	0.378	ng/ul	100
9) Fluorene	15.30	166	12244	0.369	ng/ul	100
11) Pentachlorophenol	16.65	266	1929	0.361	ng/ul	98
12) Phenanthrene	17.03	178	19834	0.417	ng/ul	99
13) Anthracene	17.12	178	19132	0.426	ng/ul	99
15) Fluoranthene	19.05	202	23043	0.394	ng/ul	96
17) Pyrene	19.42	202	23760	0.495	ng/ul	98
18) Benzo(a)anthracene	21.18	228	22042	0.458	ng/ul	99
19) Chrysene	21.23	228	20898	0.442	ng/ul	98
21) Benzo(b)fluoranthene	22.72	252	20680	0.413	ng/ul	90
22) Benzo(k)fluoranthene	22.76	252	20406	0.411	ng/ul#	85
23) Benzo(a)pyrene	23.27	252	19819	0.422	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.51	276	22948	0.417	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.53	278	18306	0.412	ng/ul	93
26) Benzo(g,h,i)perylene	26.16	276	19092	0.415	ng/ul	95

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

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