

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN112018\
 Data File : BN003633.D
 Acq On : 20 Nov 2018 16:02
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.472

Quant Time: Nov 20 16:53:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 20 03:27:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	3946	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	15532	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	8376	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	20825	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	23266	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	21588	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	9510	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	24734	0.41	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.69	128	17403	0.399	ng/ul	99
5) 2-Methylnaphthalene	12.30	142	12192	0.400	ng/ul	99
7) Acenaphthylene	14.20	152	15196	0.403	ng/ul	100
8) Acenaphthene	14.54	153	13249	0.401	ng/ul	99
9) Fluorene	15.53	166	15346	0.398	ng/ul	100
11) Pentachlorophenol	16.87	266	1620	0.432	ng/ul	99
12) Phenanthrene	17.26	178	27253	0.399	ng/ul	100
13) Anthracene	17.35	178	22632	0.403	ng/ul	99
15) Fluoranthene	19.28	202	32403	0.401	ng/ul	99
17) Pyrene	19.64	202	32161	0.385	ng/ul	99
18) Benzo(a)anthracene	21.38	228	29891	0.407	ng/ul	99
19) Chrysene	21.43	228	33274	0.396	ng/ul	99
21) Benzo(b)fluoranthene	23.01	252	34988	0.392	ng/ul	100
22) Benzo(k)fluoranthene	23.05	252	34306	0.394	ng/ul	99
23) Benzo(a)pyrene	23.61	252	30604	0.396	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.06	276	34338	0.387	ng/ul	99
25) Dibenzo(a,h)anthracene	26.08	278	27555	0.386	ng/ul	99
26) Benzo(g,h,i)perylene	26.79	276	29378	0.381	ng/ul	100

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

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