

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN010320\
 Data File : BN009307.D
 Acq On : 03 Jan 2020 18:54
 Operator : JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.412

Quant Time: Jan 03 21:53:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 03 15:42:47 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	1776	0.40	ng/ul	0.00
2) Naphthalene-d8	10.25	136	6412	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.13	164	3699	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.88	188	7486	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	4892	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	4288	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.85	152	4333	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.91	212	8441	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.30	128	7863	0.404	ng/ul	99
5) 2-Methylnaphthalene	11.92	142	5507	0.375	ng/ul	100
7) Acenaphthylene	13.84	152	7333	0.377	ng/ul	99
8) Acenaphthene	14.19	153	5767	0.378	ng/ul	99
9) Fluorene	15.18	166	6575	0.365	ng/ul	100
11) Pentachlorophenol	16.54	266	863	0.432	ng/ul	99
12) Phenanthrene	16.92	178	10392	0.399	ng/ul	99
13) Anthracene	17.01	178	9380	0.402	ng/ul	99
15) Fluoranthene	18.94	202	11157	0.375	ng/ul	99
17) Pyrene	19.31	202	11036	0.442	ng/ul	99
18) Benzo(a)anthracene	21.07	228	8772	0.404	ng/ul	99
19) Chrysene	21.12	228	8919	0.409	ng/ul	99
21) Benzo(b)fluoranthene	22.58	252	7886	0.394	ng/ul	94
22) Benzo(k)fluoranthene	22.62	252	8100	0.401	ng/ul	94
23) Benzo(a)pyrene	23.12	252	7248	0.413	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.30	276	8392	0.439	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.31	278	6841	0.450	ng/ul	97
26) Benzo(g,h,i)perylene	25.93	276	7315	0.454	ng/ul	98

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

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