

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN010320\
 Data File : BN009299.D
 Acq On : 03 Jan 2020 13:29
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.411

Quant Time: Jan 03 15:47:18 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	1587	0.40	ng/ul	0.00
2) Naphthalene-d8	10.25	136	5767	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.13	164	3467	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.88	188	7540	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	5988	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	5035	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.85	152	3973	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.91	212	9351	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.30	128	7059	0.404	ng/ul	99
5) 2-Methylnaphthalene	11.92	142	5039	0.381	ng/ul	99
7) Acenaphthylene	13.84	152	6977	0.383	ng/ul	98
8) Acenaphthene	14.19	153	5440	0.381	ng/ul	100
9) Fluorene	15.19	166	6310	0.374	ng/ul	100
11) Pentachlorophenol	16.54	266	1020	0.507	ng/ul	98
12) Phenanthrene	16.92	178	10488	0.400	ng/ul	98
13) Anthracene	17.01	178	9714	0.413	ng/ul	100
15) Fluoranthene	18.94	202	12307	0.410	ng/ul	99
17) Pyrene	19.31	202	12441	0.407	ng/ul	100
18) Benzo(a)anthracene	21.07	228	10934	0.411	ng/ul	99
19) Chrysene	21.12	228	11012	0.412	ng/ul	99
21) Benzo(b)fluoranthene	22.58	252	9438	0.401	ng/ul	97
22) Benzo(k)fluoranthene	22.62	252	9823	0.414	ng/ul	96
23) Benzo(a)pyrene	23.12	252	8477	0.411	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.30	276	8947	0.399	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.31	278	7286	0.408	ng/ul	97
26) Benzo(g,h,i)perylene	25.93	276	7726	0.409	ng/ul	99

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

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