

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080719\
 Data File : BN006980.D
 Acq On : 07 Aug 2019 10:42
 Operator : HP/JU
 Sample : SSTD0.120
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.120

Quant Time: Aug 07 17:27:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN080719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 17:13:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	5127	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	19801	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	12353	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.87	188	30846	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	26294	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	26520	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	3728	0.11	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	10256	0.10	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.28	128	6033	0.106	ng/ul#	89
5) 2-Methylnaphthalene	11.91	142	4530	0.106	ng/ul	100
7) Acenaphthylene	13.83	152	7583	0.106	ng/ul	97
8) Acenaphthene	14.18	153	5183	0.106	ng/ul	98
9) Fluorene	15.17	166	6198	0.104	ng/ul	97
12) Phenanthrene	16.90	178	10393	0.106	ng/ul	98
13) Anthracene	17.00	178	10597	0.105	ng/ul	97
15) Fluoranthene	18.94	202	12579	0.103	ng/ul	98
17) Pyrene	19.30	202	13066	0.111	ng/ul	98
18) Benzo(a)anthracene	21.07	228	12196	0.106	ng/ul	99
19) Chrysene	21.12	228	11416	0.106	ng/ul	99
21) Benzo(b)fluoranthene	22.59	252	11317	0.105	ng/ul	88
22) Benzo(k)fluoranthene	22.63	252	11103	0.107	ng/ul#	89
23) Benzo(a)pyrene	23.12	252	10659	0.105	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.31	276	11170	0.102	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.33	278	8952	0.103	ng/ul#	87
26) Benzo(g,h,i)perylene	25.95	276	9387	0.103	ng/ul#	87

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

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