

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN021119\
 Data File : BN004370.D
 Acq On : 11 Feb 2019 09:47
 Operator : JU/SJ
 Sample : SST00.102
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.102

Quant Time: Feb 11 12:13:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 11:59:16 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	969	0.40	ng/ul	0.00
2) Naphthalene-d8	10.54	136	3750	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.39	164	2171	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.14	188	6187	0.40	ng/ul	0.00
16) Chrysene-d12	21.33	240	8176	0.40	ng/ul	0.00
20) Perylene-d12	23.61	264	9051	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.13	152	580	0.10	ng/ul	0.00
14) Fluoranthene-d10	19.17	212	2009	0.11	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.59	128	1106	0.105	ng/ul#	87
5) 2-Methylnaphthalene	12.20	142	769	0.105	ng/ul	100
7) Acenaphthylene	14.11	152	1006	0.093	ng/ul#	90
8) Acenaphthene	14.45	153	860	0.099	ng/ul	95
9) Fluorene	15.44	166	1042	0.105	ng/ul#	98
12) Phenanthrene	17.18	178	2013	0.100	ng/ul	94
13) Anthracene	17.27	178	1728	0.095	ng/ul	94
15) Fluoranthene	19.20	202	2529	0.104	ng/ul	95
17) Pyrene	19.56	202	2637	0.094	ng/ul#	93
18) Benzo(a)anthracene	21.32	228	2803	0.101	ng/ul	97
19) Chrysene	21.37	228	3255	0.109	ng/ul	97
21) Benzo(b)fluoranthene	22.92	252	3739	0.109	ng/ul	85
22) Benzo(k)fluoranthene	22.96	252	3647	0.100	ng/ul#	83
23) Benzo(a)pyrene	23.51	252	3332	0.102	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	25.89	276	4329	0.128	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.91	278	3481	0.125	ng/ul#	89
26) Benzo(a,h,i)perylene	26.60	276	4022	0.141	ng/ul	93

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

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