

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112020\
 Data File : BN012655.D
 Acq On : 20 Nov 2020 11:50
 Operator : CG/JU
 Sample : SSTDO.105
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDO.105

Quant Time: Nov 20 13:46:58 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 20 13:42:54 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	836	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	3073	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	1669	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	3317	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	2961	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	2800	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	402	0.09	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	810	0.09	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.27	128	893	0.101	ng/ul#	83
5) 2-Methylnaphthalene	11.90	142	554	0.098	ng/ul	100
7) Acenaphthylene	13.82	152	768	0.106	ng/ul#	86
8) Acenaphthene	14.17	153	660	0.108	ng/ul	95
9) Fluorene	15.17	166	707	0.104	ng/ul#	91
12) Phenanthrene	16.90	178	1118	0.106	ng/ul#	88
13) Anthracene	17.00	178	916	0.101	ng/ul#	86
15) Fluoranthene	18.93	202	1293	0.105	ng/ul	88
17) Pyrene	19.30	202	1379	0.103	ng/ul#	87
18) Benzo(a)anthracene	21.06	228	1060	0.097	ng/ul	94
19) Chrysene	21.11	228	1298	0.097	ng/ul	94
21) Benzo(b)fluoranthene	22.57	252	1297	0.115	ng/ul#	61
22) Benzo(k)fluoranthene	22.61	252	1430	0.114	ng/ul#	61
23) Benzo(a)pyrene	23.11	252	1167	0.106	ng/ul#	52
24) Indeno(1,2,3-cd)pyrene	25.28	276	1461	0.103	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.29	278	1144	0.101	ng/ul#	60
26) Benzo(a,h,i)perylene	25.91	276	1312	0.101	ng/ul#	84

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

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