

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110920\
 Data File : BM027750.D
 Acq On : 09 Nov 2020 10:39
 Operator : JU/CG
 Sample : SSTD0.278
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.278

Quant Time: Nov 10 06:21:26 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM02.4-EPA-SIM-BM110920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 06:19:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	418	0.40	ng/ul	0.00
2) Naphthalene-d8	11.23	136	1378	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	818	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.72	188	1780	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	1422	0.40	ng/ul	0.00
20) Perylene-d12	24.26	264	1221	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	447	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	928	0.19	ng/ul	0.00

Target Compounds						Ovalue
3) Naphthalene	11.28	128	821	0.204	ng/ul#	89
5) 2-Methylnaphthalene	12.87	142	552	0.188	ng/ul	98
7) Acenaphthylene	14.73	152	754	0.181	ng/ul#	87
8) Acenaphthene	15.07	153	571	0.192	ng/ul	96
9) Fluorene	16.04	166	673	0.183	ng/ul#	95
11) Pentachlorophenol	17.36	266	118	0.244	ng/ul	96
12) Phenanthrene	17.76	178	1068	0.194	ng/ul	94
13) Anthracene	17.85	178	1003	0.187	ng/ul	93
15) Fluoranthene	19.73	202	1171	0.177	ng/ul	92
17) Pyrene	20.09	202	1167	0.180	ng/ul#	92
18) Benzo(a)anthracene	21.80	228	1031	0.185	ng/ul	94
19) Chrysene	21.85	228	1072	0.186	ng/ul	95
21) Benzo(b)fluoranthene	23.51	252	985	0.200	ng/ul	87
22) Benzo(k)fluoranthene	23.56	252	969	0.189	ng/ul#	87
23) Benzo(a)pyrene	24.15	252	888	0.202	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.79	276	1008	0.178	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.80	278	826	0.176	ng/ul#	85
26) Benzo(g,h,i)perylene	27.58	276	861	0.180	ng/ul#	87

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

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