

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022420\
 Data File : BM025028.D
 Acq On : 24 Feb 2020 18:59
 Operator : CG/JU
 Sample : SST00.177
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST00.177

Quant Time: Feb 25 01:04:00 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 25 00:46:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	1913	0.40	ng/ul	0.00
2) Naphthalene-d8	10.47	136	7391	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	4952	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	12141	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13369	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	14834	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	1389	0.11	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	4141	0.12	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.52	128	2323	0.110	ng/ul#	54
5) 2-Methylnaphthalene	12.14	142	1616	0.112	ng/ul	99
7) Acenaphthylene	14.04	152	2833	0.124	ng/ul#	78
8) Acenaphthene	14.39	153	1984	0.103	ng/ul	94
9) Fluorene	15.38	166	2531	0.115	ng/ul#	90
12) Phenanthrene	17.11	178	4881	0.126	ng/ul#	87
13) Anthracene	17.20	178	4161	0.125	ng/ul#	84
15) Fluoranthene	19.13	202	8001	0.192	ng/ul	90
17) Pyrene	19.49	202	7317	0.121	ng/ul#	88
18) Benzo(a)anthracene	21.24	228	6423	0.126	ng/ul#	93
19) Chrysene	21.29	228	6517	0.121	ng/ul#	92
21) Benzo(b)fluoranthene	22.80	252	5831	0.090	ng/ul#	67
22) Benzo(k)fluoranthene	22.85	252	5802	0.088	ng/ul#	64
23) Benzo(a)pyrene	23.36	252	5317	0.094	ng/ul#	56
24) Indeno(1,2,3-cd)pyrene	25.65	276	6506	0.095	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.68	278	5543	0.100	ng/ul#	72
26) Benzo(a,h,i)perylene	26.32	276	5546	0.101	ng/ul#	70

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

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