

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110920\
 Data File : BM027749.D
 Acq On : 09 Nov 2020 10:03
 Operator : JU/CG
 Sample : SST00.177
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST00.177

Quant Time: Nov 10 06:20:53 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	547	0.40	ng/ul	0.00
2) Naphthalene-d8	11.23	136	1773	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	956	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.72	188	2025	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	1571	0.40	ng/ul	0.00
20) Perylene-d12	24.26	264	1408	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	279	0.09	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	524	0.09	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.28	128	520	0.100	ng/ul	100
5) 2-Methylnaphthalene	12.86	142	351	0.093	ng/ul	100
7) Acenaphthylene	14.73	152	438	0.090	ng/ul	100
8) Acenaphthene	15.07	153	340	0.098	ng/ul	100
9) Fluorene	16.04	166	382	0.089	ng/ul	100
12) Phenanthrene	17.76	178	610	0.098	ng/ul	100
13) Anthracene	17.85	178	566	0.093	ng/ul	100
15) Fluoranthene	19.73	202	671	0.089	ng/ul	100
17) Pyrene	20.09	202	667	0.093	ng/ul	100
18) Benzo(a)anthracene	21.80	228	596	0.097	ng/ul	100
19) Chrysene	21.85	228	607	0.095	ng/ul	100
21) Benzo(b)fluoranthene	23.51	252	590	0.104	ng/ul	100
22) Benzo(k)fluoranthene	23.56	252	553	0.094	ng/ul	100
23) Benzo(a)pyrene	24.15	252	495	0.097	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.79	276	588	0.090	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.81	278	481	0.089	ng/ul	100
26) Benzo(a,h,i)perylene	27.57	276	527	0.095	ng/ul	100

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

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