

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025438.D
 Acq On : 10 Mar 2020 06:18
 Operator : CG/JU
 Sample : L1612-11
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDL3

Manual Integrations
 APPROVED

mohammad
 3/11/2020 8:00:41 AM

Quant Time: Mar 10 11:50:11 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 09 13:32:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	3314	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	14268	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9958	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	21839	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	16080	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	14689	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	8242	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	20786	0.35	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	1773	0.044	ng/ul#	91
12) Phenanthrene	17.08	178	2077m	0.030	ng/ul	
13) Anthracene	17.16	178	6879	0.114	ng/ul	97
15) Fluoranthene	19.09	202	14937	0.199	ng/ul	97
17) Pyrene	19.45	202	24552	0.340	ng/ul	98
18) Benzo(a)anthracene	21.20	228	10114	0.176	ng/ul	96
19) Chrysene	21.26	228	8508	0.132	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	28772m	0.495	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	10149m	0.167	ng/ul	
23) Benzo(a)pyrene	23.31	252	8688	0.180	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.58	276	6853	0.105	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.58	278	1921	0.035	ng/ul#	57
26) Benzo(g,h,i)perylene	26.24	276	3937	0.069	ng/ul#	90

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

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