

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025440.D
 Acq On : 10 Mar 2020 07:30
 Operator : CG/JU
 Sample : L1612-21
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDM3

Manual Integrations
 APPROVED

mohammad
 3/11/2020 8:00:53 AM

Quant Time: Mar 10 11:52:54 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	3383	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	14598	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9734	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	19892	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	14226	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	14166	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	7079	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	15059	0.28	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	2783	0.064	ng/ul#	95
5) 2-Methylnaphthalene	12.10	142	2720	0.088	ng/ul	99
7) Acenaphthylene	14.01	152	7347	0.185	ng/ul	98
8) Acenaphthene	14.35	153	1141	0.033	ng/ul	98
9) Fluorene	15.34	166	3369	0.079	ng/ul	99
12) Phenanthrene	17.07	178	20149	0.320	ng/ul	99
13) Anthracene	17.16	178	45941	0.837	ng/ul	99
15) Fluoranthene	19.09	202	26423	0.387	ng/ul	97
17) Pyrene	19.45	202	56452	0.884	ng/ul	98
18) Benzo(a)anthracene	21.20	228	31513	0.620	ng/ul	96
19) Chrysene	21.26	228	26177	0.458	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	145704m	2.597	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	49570m	0.843	ng/ul	
23) Benzo(a)pyrene	23.31	252	40222	0.863	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.58	276	46816	0.741	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.58	278	12924	0.243	ng/ul#	96
26) Benzo(a,h,i)perylene	26.24	276	25049	0.457	ng/ul	98

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

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