

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040420\
 Data File : BM025915.D
 Acq On : 03 Apr 2020 16:55
 Operator : CG/JU
 Sample : L2062-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F3E01

Manual Integrations
 APPROVED

mohammad
 4/6/2020 1:40:29 PM

Quant Time: Apr 06 09:33:29 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 02 10:49:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	3817	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	16990	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	12090	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	27209m	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	26246	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	22327	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	9435	0.31	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	25161	0.30	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.92	152	1124	0.022	ng/ul#	84
13) Anthracene	17.08	178	2723	0.036	ng/ul#	92
15) Fluoranthene	19.01	202	8495	0.080	ng/ul	99
17) Pyrene	19.37	202	21595	0.230	ng/ul	97
18) Benzo(a)anthracene	21.13	228	9315	0.099	ng/ul	94
19) Chrysene	21.18	228	17600m	0.172	ng/ul	
21) Benzo(b)fluoranthene	22.65	252	35561m	0.433	ng/ul	
22) Benzo(k)fluoranthene	22.69	252	13864m	0.157	ng/ul	
23) Benzo(a)pyrene	23.20	252	11851	0.161	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.41	276	7958	0.081	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.42	278	2225	0.027	ng/ul#	68
26) Benzo(g,h,i)perylene	26.05	276	5751	0.069	ng/ul#	92

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

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