

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021821\
 Data File : BM028826.D
 Acq On : 19 Feb 2021 18:53
 Operator : CG/JU
 Sample : M1412-01
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBG3

Manual Integrations
 APPROVED

mohammad
 2/22/2021 2:08:07 PM

Quant Time: Feb 20 01:28:08 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM020521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 20 01:10:53 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	345	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	1428	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	803	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	1564m	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	1465	0.40	ng/ul	0.00
20) Perylene-d12	23.65	264	1452	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	451	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	973	0.24	ng/ul	0.00
Target Compounds						
12) Phenanthrene	17.31	178	125	0.029	ng/ul#	33
15) Fluoranthene	19.32	202	778	0.156	ng/ul	88
17) Pyrene	19.68	202	712m	0.134	ng/ul	
18) Benzo(a)anthracene	21.41	228	407	0.081	ng/ul#	64
19) Chrysene	21.46	228	680	0.137	ng/ul#	70
21) Benzo(b)fluoranthene	22.98	252	1242m	0.250	ng/ul	
22) Benzo(k)fluoranthene	23.03	252	433m	0.087	ng/ul	
23) Benzo(a)pyrene	23.55	252	732	0.165	ng/ul#	44
24) Indeno(1,2,3-cd)pyrene	25.91	276	709	0.123	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.92	278	158	0.033	ng/ul#	1
26) Benzo(g,h,i)perylene	26.61	276	354	0.074	ng/ul#	37

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

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