

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021021\
 Data File : BM028708.D
 Acq On : 10 Feb 2021 17:30
 Operator : CG/JU
 Sample : M1310-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0A37

Manual Integrations
 APPROVED

mohammad
 2/11/2021 9:47:11 AM

Quant Time: Feb 11 00:18:34 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 : Wed Feb 10 14:28:04 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	471	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	1848	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.55	164	1005	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.28	188	1889m	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	1689	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	1694	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	592	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	1140	0.23	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.75	128	317	0.069	ng/ul#	64
5) 2-Methylnaphthalene	12.37	142	372	0.121	ng/ul	98
12) Phenanthrene	17.32	178	1019	0.193	ng/ul	94
15) Fluoranthene	19.33	202	680	0.113	ng/ul	73
17) Pyrene	19.68	202	628m	0.102	ng/ul	
18) Benzo(a)anthracene	21.41	228	328	0.057	ng/ul#	71
19) Chrysene	21.46	228	533	0.093	ng/ul#	78
21) Benzo(b)fluoranthene	22.99	252	481m	0.083	ng/ul	
22) Benzo(k)fluoranthene	23.03	252	209m	0.036	ng/ul	
23) Benzo(a)pyrene	23.56	252	304	0.059	ng/ul#	26
24) Indeno(1,2,3-cd)pyrene	25.93	276	246	0.037	ng/ul#	100
26) Benzo(g,h,i)perylene	26.62	276	218	0.039	ng/ul#	36

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

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