

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021721\
 Data File : BM028777.D
 Acq On : 17 Feb 2021 18:15
 Operator : CG/JU
 Sample : M1379-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBGY6

Manual Integrations
 APPROVED

mohammad
 2/18/2021 1:39:08 PM

Quant Time: Feb 18 05:11:49 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 : Thu Feb 18 04:11:41 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	501	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	2065	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	1116	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	2191	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	2114	0.40	ng/ul	0.00
20) Perylene-d12	23.65	264	2180	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	394	0.13	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	1045	0.18	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	130	0.025	ng/ul#	20
7) Acenaphthylene	14.26	152	127	0.028	ng/ul#	30
11) Pentachlorophenol	16.94	266	22	0.038	ng/ul#	87
12) Phenanthrene	17.32	178	224	0.037	ng/ul#	56
15) Fluoranthene	19.32	202	536	0.077	ng/ul	67
17) Pyrene	19.68	202	546m	0.071	ng/ul	
18) Benzo(a)anthracene	21.41	228	384	0.053	ng/ul#	56
19) Chrysene	21.46	228	477m	0.066	ng/ul	
21) Benzo(b)fluoranthene	22.98	252	918m	0.123	ng/ul	
22) Benzo(k)fluoranthene	23.03	252	306m	0.041	ng/ul	
23) Benzo(a)pyrene	23.55	252	626	0.094	ng/ul#	31
24) Indeno(1,2,3-cd)pyrene	25.91	276	573	0.066	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.91	278	161	0.022	ng/ul#	1
26) Benzo(g,h,i)perylene	26.60	276	554	0.077	ng/ul#	56

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

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