

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022720\
 Data File : BM025117.D
 Acq On : 27 Feb 2020 23:27
 Operator : CG/JU
 Sample : L1582-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDK0

Manual Integrations
 APPROVED

mohammad
 2/28/2020 2:40:54 PM

Quant Time: Feb 28 05:57:48 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 25 11:03:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	7	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	160	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.32	164	483	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	772m	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	1308	0.40	ng/ul	-0.01
20) Perylene-d12	23.45	264	6m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	6167	20.52	ng/ul	-0.01
14) Fluoranthene-d10	19.08	212	18441	6.96	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.51	128	10550	21.661	ng/ul#	85
5) 2-Methylnaphthalene	12.12	142	1060	3.016	ng/ul	97
8) Acenaphthene	14.38	153	1173	0.627	ng/ul	95
9) Fluorene	15.37	166	749	0.319	ng/ul#	96
12) Phenanthrene	17.10	178	2068	0.762	ng/ul#	93
13) Anthracene	17.19	178	380	0.144	ng/ul#	55
15) Fluoranthene	19.12	202	921	0.234	ng/ul	76
17) Pyrene	19.48	202	560m	0.095	ng/ul	
18) Benzo(a)anthracene	21.22	228	250	0.044	ng/ul#	53
19) Chrysene	21.28	228	257	0.046	ng/ul#	59
21) Benzo(b)fluoranthene	22.79	252	395	16.795	ng/ul#	1
22) Benzo(k)fluoranthene	22.83	252	276	11.845	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.63	276	196	7.285	ng/ul#	100
26) Benzo(g,h,i)perylene	26.28	276	166	7.322	ng/ul#	1

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

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