

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031820\
 Data File : BM025638.D
 Acq On : 18 Mar 2020 21:10
 Operator : CG/JU
 Sample : L1786-06DL 10X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC2DL

Manual Integrations
 APPROVED

mohammad
 3/19/2020 12:43:35 PM

Quant Time: Mar 19 12:15:27 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 18 05:07:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	2529	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	10305	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	6557	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	15270	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	15693	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	16703	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	422	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	1184	0.03	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.40	128	3263	0.107	ng/ul	96
5) 2-Methylnaphthalene	12.02	142	1681	0.077	ng/ul	99
7) Acenaphthylene	13.94	152	5953	0.222	ng/ul#	86
8) Acenaphthene	14.28	153	3845	0.164	ng/ul	97
9) Fluorene	15.27	166	8373	0.292	ng/ul	100
12) Phenanthrene	17.00	178	54101	1.118	ng/ul	100
13) Anthracene	17.09	178	27739	0.659	ng/ul	99
15) Fluoranthene	19.03	202	117678	2.243	ng/ul	98
17) Pyrene	19.39	202	86900	1.234	ng/ul	99
18) Benzo(a)anthracene	21.14	228	51605	0.920	ng/ul	98
19) Chrysene	21.19	228	45593	0.724	ng/ul	98
21) Benzo(b)fluoranthene	22.67	252	56995m	0.862	ng/ul	
22) Benzo(k)fluoranthene	22.71	252	24074m	0.347	ng/ul	
23) Benzo(a)pyrene	23.21	252	33362	0.607	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.43	276	20942	0.281	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.43	278	7062	0.113	ng/ul#	85
26) Benzo(a,h,i)perylene	26.07	276	13036	0.202	ng/ul#	95

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

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