

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031820\
 Data File : BM025635.D
 Acq On : 18 Mar 2020 18:46
 Operator : CG/JU
 Sample : L1786-14
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AD0

Manual Integrations
 APPROVED

mohammad
 3/19/2020 12:35:50 PM

Quant Time: Mar 18 19:20:55 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	7.58	152	2894	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	12321	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	8079	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	18571	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	19087	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	22488	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	3944	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	10667	0.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.40	128	18015	0.494	ng/ul	99
5) 2-Methylnaphthalene	12.02	142	6794	0.261	ng/ul	100
7) Acenaphthylene	13.94	152	30112	0.912	ng/ul#	86
8) Acenaphthene	14.28	153	26995	0.932	ng/ul	100
9) Fluorene	15.27	166	44074	1.249	ng/ul	99
12) Phenanthrene	17.01	178	647642	11.004	ng/ul	99
13) Anthracene	17.10	178	173430	3.386	ng/ul	99
15) Fluoranthene	19.03	202	1602587	25.119	ng/ul	98
17) Pyrene	19.39	202	1279091	14.930	ng/ul	98
18) Benzo(a)anthracene	21.14	228	871481	12.771	ng/ul	99
19) Chrysene	21.20	228	790690	10.318	ng/ul	97
21) Benzo(b)fluoranthene	22.68	252	1059032m	11.893	ng/ul	
22) Benzo(k)fluoranthene	22.71	252	410856m	4.404	ng/ul	
23) Benzo(a)pyrene	23.22	252	737629m	9.975	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.44	276	457948	4.565	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.44	278	136033	1.612	ng/ul	99
26) Benzo(a,h,i)perylene	26.08	276	313141	3.598	ng/ul	98

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

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