

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030620\
 Data File : BM025340.D
 Acq On : 07 Mar 2020 08:41
 Operator : CG/JU
 Sample : L1619-07
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDW1

Manual Integrations
 APPROVED

mohammad
 3/9/2020 3:01:11 PM

Quant Time: Mar 07 21:03:56 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 07 19:35:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	4241	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	18992	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	15374	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	33988	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	34238	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	32329	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	2671	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	8861	0.08	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	75109	1.314	ng/ul#	81
5) 2-Methylnaphthalene	12.10	142	230868	5.546	ng/ul	100
7) Acenaphthylene	14.01	152	9097	0.126	ng/ul#	86
8) Acenaphthene	14.36	153	791458	14.174	ng/ul	97
9) Fluorene	15.35	166	744357	10.635	ng/ul#	95
12) Phenanthrene	17.08	178	3269370	29.830	ng/ul	95
13) Anthracene	17.17	178	252566	2.597	ng/ul	95
15) Fluoranthene	19.10	202	2005204	14.618	ng/ul	96
17) Pyrene	19.46	202	1376383	10.600	ng/ul	97
18) Benzo(a)anthracene	21.21	228	283609	2.254	ng/ul	98
19) Chrysene	21.26	228	247437	1.789	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	139160m	1.076	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	59476m	0.438	ng/ul	
23) Benzo(a)pyrene	23.32	252	79562	0.733	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.58	276	24120	0.178	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.59	278	6754	0.060	ng/ul#	91
26) Benzo(a,h,i)perylene	26.25	276	14982	0.130	ng/ul#	95

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

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