

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031820\
 Data File : BM025628.D
 Acq On : 18 Mar 2020 14:33
 Operator : CG/JU
 Sample : L1786-06
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC2

Manual Integrations
 APPROVED

mohammad
 3/19/2020 12:35:26 PM

Quant Time: Mar 18 15:06:29 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	2202	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	8784	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5883	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	13739	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	15267	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	16874	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	3844	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	11637	0.31	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.40	128	28462	1.095	ng/ul	98
5) 2-Methylnaphthalene	12.02	142	14931	0.803	ng/ul	99
7) Acenaphthylene	13.94	152	47805	1.989	ng/ul#	86
8) Acenaphthene	14.28	153	33814	1.604	ng/ul	98
9) Fluorene	15.27	166	76625	2.983	ng/ul	100
12) Phenanthrene	17.01	178	479993	11.024	ng/ul	100
13) Anthracene	17.10	178	262375	6.925	ng/ul	99
15) Fluoranthene	19.03	202	1097435	23.251	ng/ul	99
17) Pyrene	19.39	202	803544	11.726	ng/ul	99
18) Benzo(a)anthracene	21.14	228	522781	9.578	ng/ul	98
19) Chrysene	21.19	228	437777	7.142	ng/ul	98
21) Benzo(b)fluoranthene	22.67	252	562670m	8.421	ng/ul	
22) Benzo(k)fluoranthene	22.71	252	230801m	3.297	ng/ul	
23) Benzo(a)pyrene	23.22	252	359583m	6.481	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.43	276	207320	2.754	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.44	278	67874	1.072	ng/ul	98
26) Benzo(a,h,i)perylene	26.08	276	127460	1.952	ng/ul	98

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

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