

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031820\
 Data File : BM025629.D
 Acq On : 18 Mar 2020 15:09
 Operator : CG/JU
 Sample : L1786-07
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC3

Manual Integrations
 APPROVED

mohammad
 3/19/2020 12:35:29 PM

Quant Time: Mar 18 15:47:31 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	2094	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	8504	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5767	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	13062	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	14599	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	16219	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	3373	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	9747	0.27	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.40	128	54106	2.149	ng/ul	98
5) 2-Methylnaphthalene	12.02	142	31707	1.762	ng/ul	100
7) Acenaphthylene	13.94	152	88192	3.743	ng/ul#	86
8) Acenaphthene	14.28	153	48730	2.358	ng/ul	99
9) Fluorene	15.27	166	139989	5.559	ng/ul	100
12) Phenanthrene	17.00	178	754982	18.239	ng/ul	100
13) Anthracene	17.10	178	717322	19.913	ng/ul	99
15) Fluoranthene	19.03	202	1614370	35.976	ng/ul	97
17) Pyrene	19.39	202	1184290	18.073	ng/ul	97
18) Benzo(a)anthracene	21.14	228	789934	15.134	ng/ul	98
19) Chrysene	21.19	228	675089	11.518	ng/ul	98
21) Benzo(b)fluoranthene	22.68	252	863976m	13.452	ng/ul	
22) Benzo(k)fluoranthene	22.71	252	329427m	4.896	ng/ul	
23) Benzo(a)pyrene	23.22	252	569989m	10.687	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.43	276	329228	4.550	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.44	278	105036	1.726	ng/ul	99
26) Benzo(a,h,i)perylene	26.08	276	209399	3.336	ng/ul	98

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

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