

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031820\
 Data File : BM025633.D
 Acq On : 18 Mar 2020 17:34
 Operator : CG/JU
 Sample : L1786-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC7

Manual Integrations
 APPROVED

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

Quant Time: Mar 18 18:20:08 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	2235	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	9036	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5687	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	12885	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	13396	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	14918	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	3111	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	8209	0.23	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.40	128	31176	1.165	ng/ul	99
5) 2-Methylnaphthalene	12.02	142	7765	0.406	ng/ul	100
7) Acenaphthylene	13.94	152	11903	0.512	ng/ul#	87
8) Acenaphthene	14.28	153	28724	1.409	ng/ul	99
9) Fluorene	15.27	166	43697	1.760	ng/ul	100
12) Phenanthrene	17.01	178	484472	11.865	ng/ul	99
13) Anthracene	17.10	178	131749	3.708	ng/ul	99
15) Fluoranthene	19.02	202	743498	16.796	ng/ul	97
17) Pyrene	19.39	202	562605	9.357	ng/ul	98
18) Benzo(a)anthracene	21.14	228	394896	8.245	ng/ul	99
19) Chrysene	21.19	228	344125	6.398	ng/ul	97
21) Benzo(b)fluoranthene	22.67	252	451007m	7.635	ng/ul	
22) Benzo(k)fluoranthene	22.71	252	177754m	2.872	ng/ul	
23) Benzo(a)pyrene	23.21	252	299951m	6.115	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.43	276	184596	2.774	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.44	278	58933	1.053	ng/ul	98
26) Benzo(a,h,i)perylene	26.08	276	120523	2.088	ng/ul	98

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

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