

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031320\
 Data File : BM025560.D
 Acq On : 14 Mar 2020 03:15
 Operator : CG/JU
 Sample : L1786-04
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC0

Manual Integrations
 APPROVED

mohammad
 3/16/2020 1:52:49 PM

Quant Time: Mar 16 07:11:14 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 : Mon Mar 16 06:44:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	3762	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	16741	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	12125	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	26623m	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	19949	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	18436	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	8683	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	21330	0.29	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.47	128	9666	0.195	ng/ul	100
5) 2-Methylnaphthalene	12.09	142	2015	0.057	ng/ul	99
7) Acenaphthylene	14.00	152	2695	0.054	ng/ul#	83
8) Acenaphthene	14.35	153	7820	0.180	ng/ul	99
9) Fluorene	15.33	166	7340	0.139	ng/ul	98
12) Phenanthrene	17.07	178	27993	0.332	ng/ul	100
13) Anthracene	17.16	178	9481	0.129	ng/ul	97
15) Fluoranthene	19.08	202	64151	0.701	ng/ul	97
17) Pyrene	19.45	202	55420	0.619	ng/ul	97
18) Benzo(a)anthracene	21.20	228	38145	0.535	ng/ul	98
19) Chrysene	21.25	228	35788	0.447	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	38774m	0.531	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	15975m	0.209	ng/ul	
23) Benzo(a)pyrene	23.30	252	27196	0.449	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.56	276	15451	0.188	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.57	278	4891	0.071	ng/ul#	79
26) Benzo(a,h,i)perylene	26.22	276	13395	0.188	ng/ul#	96

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

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