

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031320\
 Data File : BM025544.D
 Acq On : 13 Mar 2020 16:57
 Operator : CG/JU
 Sample : L1786-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AB7

Manual Integrations
 APPROVED

mohammad
 3/16/2020 1:52:08 PM

Quant Time: Mar 13 17:38:08 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 : Fri Mar 13 16:53:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	2800	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	12244	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	9321	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	21826m	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	16610	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	14902	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	7212	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	19460	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	10474	0.289	ng/ul	100
5) 2-Methylnaphthalene	12.09	142	2052	0.079	ng/ul	100
7) Acenaphthylene	14.00	152	4149	0.109	ng/ul#	85
8) Acenaphthene	14.35	153	12186	0.365	ng/ul	99
9) Fluorene	15.33	166	13347	0.328	ng/ul	100
12) Phenanthrene	17.07	178	40844	0.591	ng/ul	100
13) Anthracene	17.16	178	12029	0.200	ng/ul	98
15) Fluoranthene	19.08	202	64239	0.857	ng/ul	98
17) Pyrene	19.45	202	51218	0.687	ng/ul	99
18) Benzo(a)anthracene	21.20	228	25037	0.422	ng/ul	97
19) Chrysene	21.25	228	22267	0.334	ng/ul	98
21) Benzo(b)fluoranthene	22.75	252	28660m	0.486	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	11391m	0.184	ng/ul	
23) Benzo(a)pyrene	23.31	252	20039	0.409	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.57	276	11703	0.176	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.58	278	3602	0.064	ng/ul#	70
26) Benzo(a,h,i)perylene	26.23	276	10372	0.180	ng/ul#	95

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

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