

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031220\
 Data File : BM025516.D
 Acq On : 12 Mar 2020 11:05
 Operator : CG/JU
 Sample : L1677-16
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0CD8

Manual Integrations
 APPROVED

mohammad
 3/16/2020 1:49:47 PM

Quant Time: Mar 13 04:42:18 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 04:33:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	2324	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	9936	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	7084	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	16698	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	12043	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	12143	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	5464	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	14092	0.31	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	912	0.031	ng/ul#	81
5) 2-Methylnaphthalene	12.09	142	644	0.031	ng/ul	98
7) Acenaphthylene	14.00	152	710	0.025	ng/ul#	75
9) Fluorene	15.33	166	937	0.030	ng/ul#	94
12) Phenanthrene	17.07	178	11468	0.217	ng/ul	98
13) Anthracene	17.16	178	2874	0.062	ng/ul#	92
15) Fluoranthene	19.09	202	16740	0.292	ng/ul	98
17) Pyrene	19.45	202	12863m	0.238	ng/ul	
18) Benzo(a)anthracene	21.20	228	6155	0.143	ng/ul#	92
19) Chrysene	21.25	228	7188m	0.149	ng/ul	
21) Benzo(b)fluoranthene	22.75	252	7366m	0.153	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	2785m	0.055	ng/ul	
23) Benzo(a)pyrene	23.31	252	4184	0.105	ng/ul#	71
24) Indeno(1,2,3-cd)pyrene	25.57	276	3035	0.056	ng/ul#	83
25) Dibenzo(a,h)anthracene	25.59	278	1108	0.024	ng/ul#	27
26) Benzo(a,h,i)perylene	26.24	276	2698	0.057	ng/ul#	83

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

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