

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031017\
 Data File : BM009185.D
 Acq On : 10 Mar 2017 21:09
 Operator : SJ/MA
 Sample : I2107-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AJ6

Manual Integrations
 APPROVED

mohammad
 3/13/2017 2:14:16 PM

Quant Time: Mar 11 02:28:21 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 11 00:54:34 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	349	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	1141	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	854	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	1777m	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1704	0.40	ng/ul	0.00
20) Perylene-d12	23.77	264	1330m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	349	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	918m	0.17	ng/ul	0.00
Target Compounds						
						Qvalue
3) Naphthalene	10.73	128	28049	8.53	ng/ul#	93
5) 2-Methylnaphthalene	12.34	142	29125	12.31	ng/ul	100
7) Acenaphthylene	14.23	152	3178	0.72	ng/ul#	79
8) Acenaphthene	14.59	153	1213	0.36	ng/ul#	70
9) Fluorene	15.56	166	1394m	0.35	ng/ul	
12) Phenanthrene	17.30	178	40396	7.24	ng/ul	97
13) Anthracene	17.40	178	5541	1.01	ng/ul	94
15) Fluoranthene	19.33	202	69314	9.52	ng/ul	93
17) Pyrene	19.69	202	62933	10.74	ng/ul	96
18) Benzo(a)anthracene	21.42	228	33865	5.98	ng/ul	94
19) Chrysene	21.47	228	44444	8.28	ng/ul	93
21) Benzo(b)fluoranthene	23.07	252	64272m	12.30	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	19477m	4.22	ng/ul	
23) Benzo(a)pyrene	23.67	252	32097m	6.83	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.15	276	27335m	4.57	ng/ul	
25) Dibenzo(a,h)anthracene	26.16	278	7169m	1.43	ng/ul	
26) Benzo(g,h,i)perylene	26.89	276	24073m	4.72	ng/ul	

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

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