

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030917\
 Data File : BM009162.D
 Acq On : 10 Mar 2017 01:09
 Operator : SJ/MA
 Sample : I2069-06DL 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AG7DL

Manual Integrations
 APPROVED

mohammad
 3/10/2017 6:01:51 PM

Quant Time: Mar 10 04:07:15 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 10 03:16:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	258	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	784	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	423	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	965	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1063	0.40	ng/ul	0.00
20) Perylene-d12	23.78	264	1053	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	77	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	205	0.07	ng/ul	0.00

Target Compounds				Qvalue		
8) Acenaphthene	14.57	153	84	0.05	ng/ul#	83
9) Fluorene	15.56	166	181	0.09	ng/ul#	76
12) Phenanthrene	17.30	178	1764	0.58	ng/ul	94
13) Anthracene	17.40	178	530m	0.18	ng/ul	
15) Fluoranthene	19.32	202	2378	0.60	ng/ul	98
17) Pyrene	19.68	202	2081	0.57	ng/ul	97
18) Benzo(a)anthracene	21.42	228	1082	0.31	ng/ul	93
19) Chrysene	21.47	228	938m	0.28	ng/ul	
21) Benzo(b)fluoranthene	23.07	252	1372m	0.33	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	351m	0.10	ng/ul	
23) Benzo(a)pyrene	23.67	252	894	0.24	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.15	276	556	0.12	ng/ul#	87
26) Benzo(g,h,i)perylene	26.89	276	494	0.12	ng/ul#	73

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

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