

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080818\
 Data File : BE097231.D
 Acq On : 8 Aug 2018 12:32
 Operator : SJ/JU
 Sample : J4268-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC5

Manual Integrations
 APPROVED

Sohil
 8/9/2018 4:42:56 PM

Quant Time: Aug 09 00:06:09 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 23:25:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1112	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	5984	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3658	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	8651m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	8982m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	8809m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2046	0.21	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	6128m	0.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	4334	0.299	ng/ul#	94
5) 2-Methylnaphthalene	11.80	142	1999	0.209	ng/ul	99
7) Acenaphthylene	13.73	152	8922	0.564	ng/ul	98
8) Acenaphthene	14.08	153	4619	0.404	ng/ul#	95
9) Fluorene	15.08	166	6991	0.532	ng/ul	95
12) Phenanthrene	16.81	178	251458m	11.402	ng/ul	
13) Anthracene	16.90	178	31653m	1.539	ng/ul	
15) Fluoranthene	18.84	202	774515m	26.569	ng/ul	
17) Pyrene	19.21	202	666008	27.672	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	306171m	11.929	ng/ul	
19) Chrysene	21.02	228	413747	15.644	ng/ul	99
21) Benzo(b)fluoranthene	22.54	252	556599m	23.183	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	195360m	7.547	ng/ul	
23) Benzo(a)pyrene	23.12	252	361072m	15.029	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.51	276	289097m	9.615	ng/ul	
25) Dibenzo(a,h)anthracene	25.52	278	64970m	2.641	ng/ul	
26) Benzo(a,h,i)perylene	26.22	276	290924m	11.196	ng/ul	

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

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