

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080318\
 Data File : BE097176.D
 Acq On : 3 Aug 2018 11:47
 Operator : SJ/JU
 Sample : J4172-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0D00

Manual Integrations
 APPROVED

Sohil
 8/6/2018 4:32:21 PM

Quant Time: Aug 03 15:38:36 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 : Fri Aug 03 11:26:05 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1475	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7846	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4958	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	11734m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	12707m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	11522m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	3397	0.27	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	9853m	0.24	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	1059	0.056	ng/ul#	86
5) 2-Methylnaphthalene	11.81	142	389	0.031	ng/ul	98
8) Acenaphthene	14.09	153	960	0.062	ng/ul#	93
9) Fluorene	15.08	166	1095	0.061	ng/ul#	92
12) Phenanthrene	16.81	178	7596m	0.254	ng/ul	
13) Anthracene	16.90	178	2549m	0.091	ng/ul	
15) Fluoranthene	18.84	202	7963m	0.201	ng/ul	
17) Pyrene	19.21	202	7728m	0.227	ng/ul	
18) Benzo(a)anthracene	20.96	228	4248m	0.117	ng/ul	
19) Chrysene	21.01	228	3418m	0.091	ng/ul	
21) Benzo(b)fluoranthene	22.54	252	4400m	0.140	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	1950m	0.058	ng/ul	
23) Benzo(a)pyrene	23.12	252	3617	0.115	ng/ul#	75
24) Indeno(1,2,3-cd)pyrene	25.51	276	2230	0.057	ng/ul#	86
26) Benzo(g,h,i)perylene	26.20	276	1924	0.057	ng/ul#	50

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

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