

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080318\
 Data File : BE097178.D
 Acq On : 3 Aug 2018 13:01
 Operator : SJ/JU
 Sample : J4172-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0D05

Manual Integrations
 APPROVED

Sohil
 8/6/2018 4:32:22 PM

Quant Time: Aug 03 15:48:16 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	1462	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7691	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	5098	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	12037m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	12578m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	10800m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	4648	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	13918m	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	1227	0.066	ng/ul#	87
5) 2-Methylnaphthalene	11.81	142	323	0.026	ng/ul#	1
8) Acenaphthene	14.09	153	560	0.035	ng/ul#	93
9) Fluorene	15.08	166	989	0.054	ng/ul#	84
12) Phenanthrene	16.81	178	2699m	0.088	ng/ul	
15) Fluoranthene	18.84	202	2217m	0.055	ng/ul	
17) Pyrene	19.21	202	2593	0.077	ng/ul	97
18) Benzo(a)anthracene	20.96	228	1012m	0.028	ng/ul	
19) Chrysene	21.01	228	1255m	0.034	ng/ul	
21) Benzo(b)fluoranthene	22.54	252	1180	0.040	ng/ul#	1
26) Benzo(g,h,i)perylene	26.21	276	826	0.026	ng/ul#	32

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

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