

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080218\
 Data File : BE097163.D
 Acq On : 2 Aug 2018 22:36
 Operator : SJ/JU
 Sample : J4171-06
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JJ1J4

Manual Integrations
 APPROVED

Sohil
 8/3/2018 5:55:12 PM

Quant Time: Aug 03 00:23:43 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 : Thu Aug 02 14:49:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1340	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6933	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4642	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	11037m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	11942m	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	10244m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	4926	0.44	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	15286m	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	1862	0.111	ng/ul#	93
5) 2-Methylnaphthalene	11.80	142	367	0.033	ng/ul	94
12) Phenanthrene	16.81	178	3654m	0.130	ng/ul	
15) Fluoranthene	18.84	202	4534m	0.122	ng/ul	
17) Pyrene	19.21	202	5566m	0.174	ng/ul	
18) Benzo(a)anthracene	20.96	228	2318	0.068	ng/ul#	85
19) Chrysene	21.01	228	2862	0.081	ng/ul#	88
21) Benzo(b)fluoranthene	22.54	252	2867m	0.103	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	1116m	0.037	ng/ul	
23) Benzo(a)pyrene	23.11	252	2746	0.098	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.50	276	1624	0.046	ng/ul#	83
26) Benzo(g,h,i)perylene	26.20	276	1780	0.059	ng/ul#	91

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

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