

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042318\
 Data File : BE096088.D
 Acq On : 23 Apr 2018 17:01
 Operator : SJ/JU
 Sample : J2355-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Y30

Manual Integrations
 APPROVED

Sohil
 4/24/2018 7:10:01 PM

Quant Time: Apr 24 06:53:55 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 24 05:21:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	1860	0.40	ng/ul	0.00
2) Naphthalene-d8	11.20	136	5036	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2987	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6491m	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	6210m	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	7627	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	2396	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	6147m	0.23	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.70	152	668	0.043	ng/ul#	87
9) Fluorene	15.99	166	269	0.021	ng/ul#	88
12) Phenanthrene	17.72	178	4415m	0.231	ng/ul	
13) Anthracene	17.80	178	749m	0.039	ng/ul	
15) Fluoranthene	19.68	202	12808m	0.492	ng/ul	
17) Pyrene	20.04	202	13580	0.644	ng/ul#	80
18) Benzo(a)anthracene	21.75	228	8194m	0.327	ng/ul	
19) Chrysene	21.80	228	6919	0.383	ng/ul	95
21) Benzo(b)fluoranthene	23.57	252	9117m	0.369	ng/ul	
22) Benzo(k)fluoranthene	23.62	252	3281m	0.133	ng/ul	
23) Benzo(a)pyrene	24.26	252	5660	0.238	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	27.16	276	4579	0.173	ng/ul#	74
25) Dibenzo(a,h)anthracene	27.18	278	1361	0.064	ng/ul#	81
26) Benzo(g,h,i)perylene	28.04	276	3389	0.153	ng/ul	97

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

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