

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042318\
 Data File : BE096092.D
 Acq On : 23 Apr 2018 19:24
 Operator : SJ/JU
 Sample : J2355-12
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Y27

Manual Integrations
 APPROVED

Sohil
 4/24/2018 7:10:39 PM

Quant Time: Apr 24 10:47:44 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 06:59:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	2111	0.40	ng/ul	0.00
2) Naphthalene-d8	11.20	136	6080	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	3534	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	7764m	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	7288m	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	6867m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	2458	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	6537m	0.20	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.70	152	748	0.041	ng/ul#	82
8) Acenaphthene	15.03	153	359	0.027	ng/ul#	85
9) Fluorene	15.99	166	903	0.061	ng/ul	90
12) Phenanthrene	17.71	178	19117m	0.835	ng/ul	
13) Anthracene	17.80	178	1698m	0.075	ng/ul	
15) Fluoranthene	19.68	202	18050m	0.580	ng/ul	
17) Pyrene	20.04	202	30109	1.216	ng/ul#	80
18) Benzo(a)anthracene	21.75	228	11644m	0.395	ng/ul	
19) Chrysene	21.80	228	9552m	0.451	ng/ul	
21) Benzo(b)fluoranthene	23.57	252	8356m	0.376	ng/ul	
22) Benzo(k)fluoranthene	23.62	252	2102m	0.095	ng/ul	
23) Benzo(a)pyrene	24.26	252	6117m	0.285	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.16	276	3605m	0.151	ng/ul	
25) Dibenzo(a,h)anthracene	27.18	278	1153m	0.060	ng/ul	
26) Benzo(g,h,i)perylene	28.03	276	2861m	0.144	ng/ul	

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

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