

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042318\
 Data File : BE096083.D
 Acq On : 23 Apr 2018 13:59
 Operator : SJ/JU
 Sample : J2355-07MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Y33MSD

Manual Integrations
 APPROVED

Sohil
 4/24/2018 7:09:51 PM

Quant Time: Apr 24 10:34:44 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 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	1844	0.40	ng/ul	0.00
2) Naphthalene-d8	11.20	136	5081m	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2786	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6174	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	7412	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	6127	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	3054	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	8005	0.31	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.25	128	997	0.070	ng/ul#	80
5) 2-Methylnaphthalene	12.83	142	259	0.027	ng/ul	100
7) Acenaphthylene	14.70	152	1069	0.074	ng/ul#	87
8) Acenaphthene	15.03	153	3569	0.344	ng/ul	95
9) Fluorene	15.99	166	293	0.025	ng/ul#	81
11) Pentachlorophenol	17.33	266	403	0.308	ng/ul	97
12) Phenanthrene	17.71	178	5954	0.327	ng/ul#	89
13) Anthracene	17.80	178	1350	0.075	ng/ul#	90
15) Fluoranthene	19.68	202	21544	0.870	ng/ul	84
17) Pyrene	20.04	202	30281	1.202	ng/ul#	79
18) Benzo(a)anthracene	21.75	228	14066m	0.470	ng/ul	
19) Chrysene	21.80	228	10739m	0.498	ng/ul	
21) Benzo(b)fluoranthene	23.57	252	15316m	0.773	ng/ul	
22) Benzo(k)fluoranthene	23.62	252	4379m	0.222	ng/ul	
23) Benzo(a)pyrene	24.26	252	9733	0.509	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	27.17	276	7396	0.348	ng/ul#	72
25) Dibenzo(a,h)anthracene	27.17	278	1792	0.105	ng/ul#	72
26) Benzo(g,h,i)perylene	28.03	276	5814	0.327	ng/ul	98

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

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