

Data Path : Z:\HPCHEM1\BNA E\DATA\BE103117\
 Data File : BE094683.D
 Acq On : 1 Nov 2017 6:05
 Operator : SJ/JU
 Sample : I5844-20DL 5X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AH4DL

Manual Integrations
 APPROVED

Sohil
 11/1/2017 4:30:02 PM

Quant Time: Nov 01 09:56:23 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 09:37:39 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1408	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	8358	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	4621	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	10131	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	10458	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	11014	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	449	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	1087	0.04	ng/ul	-0.01
Target Compounds						Ovalue
3) Naphthalene	10.74	128	460	0.023	ng/ul#	68
5) 2-Methylnaphthalene	12.36	142	352	0.025	ng/ul	99
7) Acenaphthylene	14.24	152	1201	0.057	ng/ul#	87
8) Acenaphthene	14.57	153	1584	0.105	ng/ul	96
9) Fluorene	15.57	166	1725	0.102	ng/ul	97
12) Phenanthrene	17.29	178	128628	4.807	ng/ul#	92
13) Anthracene	17.38	178	10097	0.382	ng/ul#	93
15) Fluoranthene	19.30	202	374160	12.094	ng/ul	84
17) Pyrene	19.67	202	289334	8.949	ng/ul#	80
18) Benzo(a)anthracene	21.40	228	100223	3.352	ng/ul#	87
19) Chrysene	21.46	228	170109	5.505	ng/ul	98
21) Benzo(b)fluoranthene	23.18	252	208262m	6.722	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	90065m	2.713	ng/ul	
23) Benzo(a)pyrene	23.84	252	109807	3.546	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	26.65	276	83899	2.618	ng/ul#	76
25) Dibenzo(a,h)anthracene	26.64	278	17910	0.663	ng/ul	92
26) Benzo(a,h,i)perylene	27.49	276	65929	2.562	ng/ul	97

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

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