

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094375.D
 Acq On : 15 Oct 2017 3:23
 Operator : SJ/JU
 Sample : I5548-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB2

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:25:38 PM

Quant Time: Oct 16 05:43:44 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 05:09:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1641	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	9004	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5372	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11719	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	11228	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	10211m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	2622	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	6408	0.16	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.36	142	516	0.036	ng/ul	99
7) Acenaphthylene	14.25	152	890	0.042	ng/ul#	81
8) Acenaphthene	14.59	153	972	0.056	ng/ul	97
9) Fluorene	15.57	166	1070	0.053	ng/ul#	88
12) Phenanthrene	17.29	178	41971	1.371	ng/ul#	87
13) Anthracene	17.38	178	7909	0.266	ng/ul#	90
15) Fluoranthene	19.31	202	147722	3.671	ng/ul	84
17) Pyrene	19.67	202	125906	4.030	ng/ul#	82
18) Benzo(a)anthracene	21.41	228	73424	2.465	ng/ul#	88
19) Chrysene	21.46	228	81486	2.369	ng/ul	98
21) Benzo(b)fluoranthene	23.17	252	92826	3.347	ng/ul	88
22) Benzo(k)fluoranthene	23.21	252	36344m	1.134	ng/ul	
23) Benzo(a)pyrene	23.83	252	66991m	2.324	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.63	276	45572m	1.546	ng/ul	
25) Dibenzo(a,h)anthracene	26.63	278	12558m	0.508	ng/ul	
26) Benzo(a,h,i)perylene	27.46	276	36239m	1.425	ng/ul	

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

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