

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094394.D
 Acq On : 15 Oct 2017 15:40
 Operator : SJ/JU
 Sample : I5548-08DL 5X
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB9DL

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:26:23 PM

Quant Time: Oct 16 07:36:01 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 06:50:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1789	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	9910	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	6185	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	14316	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	14531	0.40	ng/ul	0.00
20) Perylene-d12	23.94	264	14161	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	989	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	2807	0.06	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.75	128	1415	0.058	ng/ul#	85
5) 2-Methylnaphthalene	12.35	142	984	0.062	ng/ul	99
7) Acenaphthylene	14.24	152	952	0.039	ng/ul#	83
8) Acenaphthene	14.58	153	11401	0.574	ng/ul	98
9) Fluorene	15.57	166	13474	0.580	ng/ul	95
12) Phenanthrene	17.29	178	234237	6.261	ng/ul#	91
13) Anthracene	17.38	178	50383	1.385	ng/ul#	90
15) Fluoranthene	19.31	202	325167	6.615	ng/ul	84
17) Pyrene	19.67	202	257092	6.359	ng/ul#	82
18) Benzo(a)anthracene	21.40	228	130549	3.387	ng/ul#	86
19) Chrysene	21.45	228	114101	2.563	ng/ul	99
21) Benzo(b)fluoranthene	23.17	252	129298m	3.361	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	54692m	1.230	ng/ul	
23) Benzo(a)pyrene	23.82	252	92678	2.318	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.61	276	56234	1.375	ng/ul#	79
25) Dibenzo(a,h)anthracene	26.61	278	14972	0.437	ng/ul	95
26) Benzo(a,h,i)perylene	27.44	276	44607	1.265	ng/ul	99

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

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