

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094641.D
 Acq On : 30 Oct 2017 10:52
 Operator : SJ/JU
 Sample : I5842-20
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB9

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:40:31 PM

Quant Time: Oct 30 18:38:29 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 30 04:29:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1239	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	7506	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	4289	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8827	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	8568	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	7415m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	2700	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	6032	0.23	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.24	152	2878	0.148	ng/ul	98
8) Acenaphthene	14.58	153	526	0.038	ng/ul#	93
9) Fluorene	15.57	166	602	0.038	ng/ul#	85
12) Phenanthrene	17.29	178	15756m	0.676	ng/ul	
13) Anthracene	17.39	178	9849	0.428	ng/ul	95
15) Fluoranthene	19.31	202	67400	2.500	ng/ul	84
17) Pyrene	19.68	202	120426	4.547	ng/ul#	79
18) Benzo(a)anthracene	21.41	228	60270	2.461	ng/ul#	88
19) Chrysene	21.47	228	70671	2.791	ng/ul	97
21) Benzo(b)fluoranthene	23.19	252	91914m	4.407	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	38493m	1.722	ng/ul	
23) Benzo(a)pyrene	23.86	252	41678m	1.999	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.68	276	14416m	0.668	ng/ul	
25) Dibenzo(a,h)anthracene	26.66	278	5232m	0.288	ng/ul	
26) Benzo(g,h,i)perylene	27.52	276	8849m	0.511	ng/ul	

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

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