

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094720.D
 Acq On : 3 Nov 2017 22:42
 Operator : SJ/JU
 Sample : I5902-05 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-65(0.5-1.5)-101317

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:16:45 PM

Quant Time: Nov 04 00:35:59 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 04 01:15:30 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1366	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	7682	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	4367	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9510	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	9276	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	9521	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	498	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	1187	0.04	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	965	0.051	ng/ul#	80
5) 2-Methylnaphthalene	12.35	142	850	0.065	ng/ul	97
7) Acenaphthylene	14.24	152	1014	0.051	ng/ul#	85
8) Acenaphthene	14.58	153	301	0.021	ng/ul#	87
9) Fluorene	15.56	166	536	0.034	ng/ul#	80
12) Phenanthrene	17.29	178	11261	0.448	ng/ul#	94
13) Anthracene	17.38	178	1347m	0.054	ng/ul	
15) Fluoranthene	19.30	202	19923	0.686	ng/ul	84
17) Pyrene	19.67	202	22807	0.795	ng/ul#	81
18) Benzo(a)anthracene	21.40	228	10177	0.384	ng/ul#	83
19) Chrysene	21.46	228	15636	0.570	ng/ul#	84
21) Benzo(b)fluoranthene	23.18	252	20105m	0.751	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	5813m	0.203	ng/ul	
23) Benzo(a)pyrene	23.84	252	11045	0.413	ng/ul#	51
24) Indeno(1,2,3-cd)pyrene	26.66	276	8466	0.306	ng/ul#	74
25) Dibenzo(a,h)anthracene	26.65	278	2054	0.088	ng/ul#	1
26) Benzo(g,h,i)perylene	27.51	276	7965	0.358	ng/ul#	66

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

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