

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094735.D
 Acq On : 4 Nov 2017 7:33
 Operator : SJ/JU
 Sample : I5955-02 5X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-10(3-5)-101617

Manual Integrations
 APPROVED

Sohil
 11/6/2017 5:58:40 PM

Quant Time: Nov 06 00:08:24 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 06:57:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1165	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.69	136	7198	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.51	164	4208	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9825	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	10338	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	12288	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	1010	0.09	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	2511	0.09	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.24	152	489	0.026	ng/ul#	65
8) Acenaphthene	14.57	153	717	0.052	ng/ul	96
9) Fluorene	15.56	166	688	0.045	ng/ul#	89
12) Phenanthrene	17.29	178	8866	0.342	ng/ul#	92
13) Anthracene	17.38	178	2010	0.078	ng/ul	96
15) Fluoranthene	19.31	202	14402	0.480	ng/ul	84
17) Pyrene	19.67	202	12982	0.406	ng/ul#	83
18) Benzo(a)anthracene	21.41	228	6855	0.232	ng/ul#	88
19) Chrysene	21.46	228	6329m	0.207	ng/ul	
21) Benzo(b)fluoranthene	23.18	252	7235m	0.209	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	2450m	0.066	ng/ul	
23) Benzo(a)pyrene	23.85	252	4892	0.142	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	26.67	276	2630	0.074	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.66	278	675	0.022	ng/ul#	6
26) Benzo(g,h,i)perylene	27.52	276	2179	0.076	ng/ul#	65

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

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