

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101317\
 Data File : BE094346.D
 Acq On : 14 Oct 2017 5:11
 Operator : SJ/JU
 Sample : I5649-01DL 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-39(0.5-1.5)-100317DL

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:23:20 PM

Quant Time: Oct 14 22:18:21 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 14 05:22:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1633	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	8441	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	4685	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9883	0.40	ng/ul	-0.02
16) Chrysene-d12	21.44	240	9908	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	10024	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	1096	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	2791	0.08	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	2514	0.12	ng/ul#	95
5) 2-Methylnaphthalene	12.37	142	716	0.05	ng/ul	98
7) Acenaphthylene	14.25	152	1311	0.07	ng/ul#	82
8) Acenaphthene	14.59	153	1007	0.07	ng/ul	97
9) Fluorene	15.58	166	1031	0.06	ng/ul#	91
12) Phenanthrene	17.31	178	31383	1.22	ng/ul#	93
13) Anthracene	17.40	178	7537	0.30	ng/ul	94
15) Fluoranthene	19.32	202	61403	1.81	ng/ul	84
17) Pyrene	19.68	202	66480	2.41	ng/ul#	84
18) Benzo(a)anthracene	21.42	228	35573	1.35	ng/ul#	89
19) Chrysene	21.47	228	33830	1.11	ng/ul	93
21) Benzo(b)fluoranthene	23.19	252	38671m	1.42	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	13288m	0.42	ng/ul	
23) Benzo(a)pyrene	23.85	252	33272	1.18	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	26.66	276	20024	0.69	ng/ul#	77
25) Dibenzo(a,h)anthracene	26.66	278	4796	0.20	ng/ul#	30
26) Benzo(g,h,i)perylene	27.49	276	18573	0.74	ng/ul#	84

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

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