

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094738.D
 Acq On : 4 Nov 2017 9:18
 Operator : SJ/JU
 Sample : I5955-05 5X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-15(0.25-1.25)-101717

Manual Integrations
 APPROVED

Sohil
 11/6/2017 5:58:47 PM

Quant Time: Nov 06 00:22:32 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	1927	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.69	136	11339	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.51	164	6323	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	13484	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	13356	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	14543	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	729	0.04	ng/ul	-0.02
14) Fluoranthene-d10	19.28	212	1663	0.04	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.73	128	5378	0.194	ng/ul#	95
5) 2-Methylnaphthalene	12.35	142	2880	0.148	ng/ul	96
7) Acenaphthylene	14.23	152	3480	0.122	ng/ul	97
12) Phenanthrene	17.29	178	8060	0.226	ng/ul	93
13) Anthracene	17.38	178	2272m	0.065	ng/ul	
15) Fluoranthene	19.31	202	9182	0.223	ng/ul	83
17) Pyrene	19.67	202	8821	0.214	ng/ul#	84
18) Benzo(a)anthracene	21.41	228	5579	0.146	ng/ul#	56
19) Chrysene	21.46	228	6968	0.177	ng/ul#	67
21) Benzo(b)fluoranthene	23.18	252	22150m	0.541	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	5650m	0.129	ng/ul	
23) Benzo(a)pyrene	23.85	252	5942	0.145	ng/ul#	53
24) Indeno(1,2,3-cd)pyrene	26.66	276	13288	0.314	ng/ul#	76
25) Dibenzo(a,h)anthracene	26.65	278	3762	0.106	ng/ul#	24
26) Benzo(g,h,i)perylene	27.52	276	12159	0.358	ng/ul#	78

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

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