

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094739.D
 Acq On : 4 Nov 2017 9:53
 Operator : SJ/JU
 Sample : I5955-06MS 5X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 11/6/2017 5:58:50 PM

Quant Time: Nov 06 00:24:56 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	1669	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.69	136	9690	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.51	164	5539	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12045	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	12004	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	11923	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	418	0.03	ng/ul	-0.02
14) Fluoranthene-d10	19.28	212	1037	0.03	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	2463	0.104	ng/ul#	92
5) 2-Methylnaphthalene	12.35	142	1706	0.103	ng/ul	98
7) Acenaphthylene	14.24	152	690	0.028	ng/ul#	68
8) Acenaphthene	14.57	153	587	0.033	ng/ul#	93
12) Phenanthrene	17.29	178	6101	0.192	ng/ul#	93
13) Anthracene	17.38	178	1093m	0.035	ng/ul	
15) Fluoranthene	19.31	202	10854	0.295	ng/ul	83
17) Pyrene	19.67	202	15153	0.408	ng/ul#	84
18) Benzo(a)anthracene	21.41	228	5592	0.163	ng/ul#	30
19) Chrysene	21.46	228	14626	0.412	ng/ul#	53
21) Benzo(b)fluoranthene	23.19	252	11257m	0.336	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	2968m	0.083	ng/ul	
23) Benzo(a)pyrene	23.85	252	6232	0.186	ng/ul#	26
24) Indeno(1,2,3-cd)pyrene	26.67	276	4777	0.138	ng/ul#	72
25) Dibenzo(a,h)anthracene	26.67	278	1445	0.049	ng/ul#	1
26) Benzo(g,h,i)perylene	27.52	276	5622	0.202	ng/ul#	50

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

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