

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062317\
 Data File : BE093338.D
 Acq On : 23 Jun 2017 15:09
 Operator : SJ/JU
 Sample : I3625-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A82

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:32:50 AM

Quant Time: Jun 24 04:26:04 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 24 03:58:33 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1603	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	8756	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5444	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	12964	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13745	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	10467m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	2014	0.16	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	5441	0.15	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	3183	0.147	ng/ul#	94
5) 2-Methylnaphthalene	12.18	142	769	0.052	ng/ul	100
7) Acenaphthylene	14.08	152	29819	1.276	ng/ul	96
8) Acenaphthene	14.42	153	544	0.029	ng/ul#	89
9) Fluorene	15.41	166	1005	0.045	ng/ul#	91
12) Phenanthrene	17.13	178	15841	0.444	ng/ul#	92
13) Anthracene	17.22	178	27660	0.891	ng/ul#	90
15) Fluoranthene	19.15	202	92635	2.118	ng/ul	84
17) Pyrene	19.51	202	161355	3.981	ng/ul#	83
18) Benzo(a)anthracene	21.24	228	103120	2.813	ng/ul#	84
19) Chrysene	21.29	228	137363	3.612	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	330395m	10.401	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	116118m	3.442	ng/ul	
23) Benzo(a)pyrene	23.55	252	163873	5.456	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.18	276	140265	4.033	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.19	278	36504m	1.246	ng/ul	
26) Benzo(g,h,i)perylene	26.95	276	103586	3.473	ng/ul	97

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

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