

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062317\
 Data File : BE093348.D
 Acq On : 23 Jun 2017 22:00
 Operator : SJ/JU
 Sample : I3627-08
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A47

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:33:28 AM

Quant Time: Jun 28 13:15:11 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 04:29:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1584	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	8358	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5364	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	12451	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	12803	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	10881m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	3608	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	9937	0.29	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.56	128	34006	1.640	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	2370	0.167	ng/ul	99
7) Acenaphthylene	14.08	152	1356	0.059	ng/ul#	88
8) Acenaphthene	14.42	153	2209	0.119	ng/ul	98
9) Fluorene	15.41	166	2800	0.129	ng/ul	92
12) Phenanthrene	17.14	178	14696	0.429	ng/ul#	91
13) Anthracene	17.22	178	7119	0.239	ng/ul#	91
15) Fluoranthene	19.15	202	12081	0.288	ng/ul	84
17) Pyrene	19.51	202	11828	0.313	ng/ul#	88
18) Benzo(a)anthracene	21.24	228	3289	0.096	ng/ul#	86
19) Chrysene	21.29	228	3286	0.093	ng/ul	91
21) Benzo(b)fluoranthene	22.93	252	3322	0.101	ng/ul#	80
22) Benzo(k)fluoranthene	22.97	252	1113m	0.032	ng/ul	
23) Benzo(a)pyrene	23.55	252	2188m	0.070	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.18	276	1173m	0.032	ng/ul	
26) Benzo(g,h,i)perylene	26.95	276	20937	0.675	ng/ul	97

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

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