

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062417\
 Data File : BE093369.D
 Acq On : 24 Jun 2017 19:35
 Operator : SJ/JU
 Sample : I3627-08DL 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A47DL

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:34:35 AM

Quant Time: Jun 26 02:04:43 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 26 01:58:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1542	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	7959	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	4807	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	11607	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	12646	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	12911	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	1776	0.15	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	4984	0.16	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	17030	0.862	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	1181	0.087	ng/ul	99
7) Acenaphthylene	14.08	152	702	0.034	ng/ul#	80
8) Acenaphthene	14.42	153	1108	0.067	ng/ul	96
9) Fluorene	15.42	166	1341	0.069	ng/ul	94
12) Phenanthrene	17.14	178	7177	0.225	ng/ul#	91
13) Anthracene	17.22	178	3630	0.131	ng/ul#	92
15) Fluoranthene	19.15	202	6193	0.158	ng/ul	84
17) Pyrene	19.51	202	6275	0.168	ng/ul#	91
18) Benzo(a)anthracene	21.24	228	1756	0.052	ng/ul#	86
19) Chrysene	21.29	228	1931	0.055	ng/ul#	89
21) Benzo(b)fluoranthene	22.93	252	1977m	0.050	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	644m	0.015	ng/ul	
23) Benzo(a)pyrene	23.55	252	1223	0.033	ng/ul#	70
24) Indeno(1,2,3-cd)pyrene	26.18	276	701	0.016	ng/ul#	76
26) Benzo(g,h,i)perylene	26.95	276	11083	0.301	ng/ul	95

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

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