

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062417\
 Data File : BE093364.D
 Acq On : 24 Jun 2017 16:30
 Operator : SJ/JU
 Sample : I3625-18DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A00DL

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:34:23 AM

Quant Time: Jun 26 01:52:06 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 01:13:30 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1236	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	6623	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	3932	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	9981	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	11304	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	9260	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	565	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	1752	0.06	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	1718	0.105	ng/ul#	93
5) 2-Methylnaphthalene	12.18	142	1018	0.091	ng/ul	99
7) Acenaphthylene	14.08	152	4364	0.259	ng/ul	99
8) Acenaphthene	14.42	153	2830	0.208	ng/ul	98
9) Fluorene	15.42	166	4104	0.257	ng/ul	94
12) Phenanthrene	17.13	178	115718	4.214	ng/ul#	90
13) Anthracene	17.22	178	21390	0.895	ng/ul#	91
15) Fluoranthene	19.15	202	220558	6.549	ng/ul	84
17) Pyrene	19.51	202	211383	6.341	ng/ul#	84
18) Benzo(a)anthracene	21.24	228	111701	3.705	ng/ul#	86
19) Chrysene	21.29	228	111109	3.552	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	122988m	4.377	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	37333m	1.251	ng/ul	
23) Benzo(a)pyrene	23.55	252	78506	2.955	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.18	276	46816	1.522	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.19	278	14326	0.553	ng/ul#	90
26) Benzo(g,h,i)perylene	26.96	276	39381	1.493	ng/ul	97

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

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