

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062417\
 Data File : BE093358.D
 Acq On : 24 Jun 2017 12:46
 Operator : SJ/JU
 Sample : I3625-11DL 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5D15DL

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:34:05 AM

Quant Time: Jun 26 01:31:16 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	1297	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	6734	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	3808	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	9698	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	11679	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	9707	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	171	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	994	0.04	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	310	0.019	ng/ul#	66
7) Acenaphthylene	14.08	152	7623	0.466	ng/ul	98
8) Acenaphthene	14.42	153	430	0.033	ng/ul#	93
9) Fluorene	15.42	166	1006	0.065	ng/ul#	89
12) Phenanthrene	17.14	178	35362	1.325	ng/ul#	92
13) Anthracene	17.22	178	11010	0.474	ng/ul#	91
15) Fluoranthene	19.15	202	173185	5.292	ng/ul	84
17) Pyrene	19.51	202	177589	5.156	ng/ul#	87
18) Benzo(a)anthracene	21.24	228	116957	3.754	ng/ul#	85
19) Chrysene	21.29	228	116966	3.619	ng/ul	99
21) Benzo(b)fluoranthene	22.93	252	156795m	5.323	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	48340m	1.545	ng/ul	
23) Benzo(a)pyrene	23.55	252	102889	3.694	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.18	276	67482	2.092	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.19	278	19517	0.719	ng/ul	93
26) Benzo(g,h,i)perylene	26.96	276	57477	2.078	ng/ul	97

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

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