

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062417\
 Data File : BE093370.D
 Acq On : 24 Jun 2017 20:14
 Operator : SJ/JU
 Sample : I3627-09DL 10X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A69DL

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:34:38 AM

Quant Time: Jun 26 02:10:23 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	1557	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	8201	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	4883	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	12232	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13844	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	12228	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	296	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	925	0.03	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.08	152	4213	0.201	ng/ul	99
12) Phenanthrene	17.14	178	3200	0.095	ng/ul#	92
13) Anthracene	17.22	178	5175m	0.177	ng/ul	
15) Fluoranthene	19.15	202	34880	0.845	ng/ul	84
17) Pyrene	19.51	202	44887	1.099	ng/ul#	87
18) Benzo(a)anthracene	21.24	228	28623	0.775	ng/ul#	85
19) Chrysene	21.29	228	31507	0.823	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	58015m	1.563	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	14927m	0.379	ng/ul	
23) Benzo(a)pyrene	23.55	252	26751	0.762	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.17	276	17778	0.438	ng/ul#	88
25) Dibenzo(a,h)anthracene	26.18	278	4796	0.140	ng/ul#	85
26) Benzo(g,h,i)perylene	26.95	276	12439	0.357	ng/ul	96

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

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