

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062417\
 Data File : BE093365.D
 Acq On : 24 Jun 2017 17:09
 Operator : SJ/JU
 Sample : I3625-08DL 10X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5B82DL

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:34:26 AM

Quant Time: Jun 26 01:54:59 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	1431	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	7651	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	4499	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	11395	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	12554	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	11527	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	424	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	1417	0.05	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	1825	0.096	ng/ul#	95
5) 2-Methylnaphthalene	12.18	142	423	0.033	ng/ul	95
7) Acenaphthylene	14.08	152	8203	0.425	ng/ul	97
12) Phenanthrene	17.14	178	5667	0.181	ng/ul#	92
13) Anthracene	17.22	178	15922m	0.583	ng/ul	
15) Fluoranthene	19.15	202	24070	0.626	ng/ul	83
17) Pyrene	19.51	202	26277	0.710	ng/ul#	85
18) Benzo(a)anthracene	21.24	228	21723	0.649	ng/ul#	83
19) Chrysene	21.29	228	25735	0.741	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	56361m	1.611	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	16373m	0.441	ng/ul	
23) Benzo(a)pyrene	23.55	252	26825	0.811	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.18	276	30146	0.787	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.19	278	7390	0.229	ng/ul#	90
26) Benzo(g,h,i)perylene	26.96	276	29441	0.896	ng/ul	96

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

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