

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062417\
 Data File : BE093368.D
 Acq On : 24 Jun 2017 18:59
 Operator : SJ/JU
 Sample : I3627-03DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A09DL

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:34:32 AM

Quant Time: Jun 26 02:02:42 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	1506	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	8021	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	4802	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	12307	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13536	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	11167	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	591	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	1744	0.05	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	6011	0.302	ng/ul#	92
5) 2-Methylnaphthalene	12.18	142	1043	0.077	ng/ul	100
7) Acenaphthylene	14.08	152	10864	0.527	ng/ul	97
8) Acenaphthene	14.42	153	1558	0.094	ng/ul	98
9) Fluorene	15.41	166	2919	0.150	ng/ul	92
12) Phenanthrene	17.14	178	22501	0.664	ng/ul#	90
13) Anthracene	17.22	178	13420	0.455	ng/ul#	90
15) Fluoranthene	19.15	202	56159	1.352	ng/ul	84
17) Pyrene	19.51	202	96531	2.418	ng/ul#	88
18) Benzo(a)anthracene	21.24	228	59747	1.655	ng/ul#	85
19) Chrysene	21.29	228	82944	2.215	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	141720m	4.182	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	47963m	1.333	ng/ul	
23) Benzo(a)pyrene	23.56	252	81539	2.545	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.18	276	52833	1.424	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.19	278	14397m	0.461	ng/ul	
26) Benzo(g,h,i)perylene	26.96	276	38875	1.222	ng/ul	96

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

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