

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061917\
 Data File : BE093266.D
 Acq On : 20 Jun 2017 00:43
 Operator : SJ/MA
 Sample : I3600-11
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A18

Manual Integrations
 APPROVED

mohammad
 6/20/2017 3:13:08 PM

Quant Time: Jun 20 09:22:03 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 20 08:49:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2576	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	13231	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	8401	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	20502m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	20271	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	15865m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	7118	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	20825	0.37	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	559	0.017	ng/ul#	71
7) Acenaphthylene	14.09	152	7035	0.195	ng/ul	97
8) Acenaphthene	14.43	153	660	0.023	ng/ul	98
12) Phenanthrene	17.14	178	1819	0.032	ng/ul#	92
13) Anthracene	17.22	178	20164	0.411	ng/ul#	87
15) Fluoranthene	19.15	202	11151	0.161	ng/ul	84
17) Pyrene	19.52	202	34925	0.584	ng/ul#	83
18) Benzo(a)anthracene	21.24	228	14214	0.263	ng/ul#	84
19) Chrysene	21.29	228	22853	0.407	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	53031	1.101	ng/ul	89
22) Benzo(k)fluoranthene	22.98	252	16634m	0.325	ng/ul	
23) Benzo(a)pyrene	23.56	252	24215m	0.532	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.18	276	16953m	0.322	ng/ul	
25) Dibenzo(a,h)anthracene	26.20	278	4726m	0.106	ng/ul	
26) Benzo(g,h,i)perylene	26.96	276	14644	0.324	ng/ul	94

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

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