

Data Path : Z:\HPCHEM1\BNA E\DATA\BE062117\
 Data File : BE093318.D
 Acq On : 21 Jun 2017 16:42
 Operator : SJ/JU
 Sample : I3627-14DL 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5E25DL

Manual Integrations
 APPROVED

mohammad
 6/22/2017 11:49:00 AM

Quant Time: Jun 22 03:36:27 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 22 02:10:17 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1652	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	9023	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5504	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13702	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13818	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	12480	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	621	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	1658	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.56	128	4214	0.19	ng/ul#	93
5) 2-Methylnaphthalene	12.18	142	485	0.03	ng/ul	98
7) Acenaphthylene	14.09	152	2750	0.12	ng/ul	99
12) Phenanthrene	17.14	178	1939	0.05	ng/ul	93
13) Anthracene	17.22	178	4773m	0.15	ng/ul	
15) Fluoranthene	19.15	202	26976	0.58	ng/ul	84
17) Pyrene	19.51	202	35530	0.87	ng/ul#	88
18) Benzo(a)anthracene	21.24	228	18055	0.49	ng/ul#	85
19) Chrysene	21.29	228	19668	0.51	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	26193m	0.69	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	11204m	0.28	ng/ul	
23) Benzo(a)pyrene	23.55	252	14032	0.39	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.17	276	7931	0.19	ng/ul#	89
25) Dibenzo(a,h)anthracene	26.19	278	2228	0.06	ng/ul#	76
26) Benzo(g,h,i)perylene	26.95	276	5139	0.14	ng/ul	94

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

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