

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061517\
 Data File : BE093219.D
 Acq On : 15 Jun 2017 21:46
 Operator : SJ/MA
 Sample : I3521-17DL 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5C60DL

Manual Integrations
 APPROVED

mohammad
 6/19/2017 11:06:50 AM

Quant Time: Jun 16 05:22:18 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 16 04:43:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2839	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	14853	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	9005	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	21017	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	21295	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	19621	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	1718	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	4304	0.08	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	1916	0.052	ng/ul#	95
5) 2-Methylnaphthalene	12.19	142	504	0.020	ng/ul	97
7) Acenaphthylene	14.09	152	7455	0.193	ng/ul	98
12) Phenanthrene	17.13	178	5600	0.097	ng/ul#	87
13) Anthracene	17.22	178	8157	0.162	ng/ul#	88
15) Fluoranthene	19.16	202	50331	0.710	ng/ul	83
17) Pyrene	19.52	202	70388	1.121	ng/ul#	82
18) Benzo(a)anthracene	21.25	228	33039	0.582	ng/ul#	85
19) Chrysene	21.30	228	40893	0.694	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	83314m	1.399	ng/ul	
22) Benzo(k)fluoranthene	22.98	252	28164m	0.445	ng/ul	
23) Benzo(a)pyrene	23.56	252	41885	0.744	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.18	276	29743	0.456	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.20	278	7181	0.131	ng/ul#	86
26) Benzo(g,h,i)perylene	26.96	276	24617	0.440	ng/ul	97

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

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