

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061417\
 Data File : BE093189.D
 Acq On : 15 Jun 2017 3:07
 Operator : SJ/MA
 Sample : I3521-16
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5C36

Manual Integrations
 APPROVED

mohammad
 6/15/2017 3:53:55 PM

Quant Time: Jun 15 04:50:56 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 15 03:59:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	4137	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	21388	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	12986	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	31453	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	27483	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	22609m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	7194	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	17584	0.21	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	18222	0.343	ng/ul#	92
5) 2-Methylnaphthalene	12.19	142	4275	0.118	ng/ul	98
7) Acenaphthylene	14.08	152	73549	1.320	ng/ul	96
8) Acenaphthene	14.43	153	3847	0.086	ng/ul	97
9) Fluorene	15.42	166	3633	0.069	ng/ul#	87
12) Phenanthrene	17.14	178	240656m	2.781	ng/ul	
13) Anthracene	17.24	178	86108	1.143	ng/ul	96
15) Fluoranthene	19.16	202	1007884	9.497	ng/ul	84
17) Pyrene	19.52	202	1044342	12.885	ng/ul#	84
18) Benzo(a)anthracene	21.24	228	504479	6.882	ng/ul#	84
19) Chrysene	21.30	228	637553	8.384	ng/ul	98
21) Benzo(b)fluoranthene	22.94	252	998157	14.548	ng/ul	88
22) Benzo(k)fluoranthene	22.98	252	336455m	4.617	ng/ul	
23) Benzo(a)pyrene	23.57	252	453412m	6.989	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.19	276	316435m	4.212	ng/ul	
25) Dibenzo(a,h)anthracene	26.21	278	82271m	1.301	ng/ul	
26) Benzo(g,h,i)perylene	26.97	276	245871	3.817	ng/ul	96

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

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