

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061417\
 Data File : BE093187.D
 Acq On : 15 Jun 2017 1:55
 Operator : SJ/MA
 Sample : I3522-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5D30

Manual Integrations
 APPROVED

mohammad
 6/15/2017 3:53:52 PM

Quant Time: Jun 15 04:44:26 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	3455	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	17586	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	10454	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	24813	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	28810	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	23939m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	4835	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	13623	0.20	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	8785	0.201	ng/ul#	94
5) 2-Methylnaphthalene	12.19	142	2172	0.073	ng/ul	99
7) Acenaphthylene	14.09	152	172155	3.837	ng/ul	96
8) Acenaphthene	14.43	153	3687	0.102	ng/ul	96
9) Fluorene	15.42	166	5286	0.125	ng/ul	91
12) Phenanthrene	17.14	178	112227	1.644	ng/ul#	83
13) Anthracene	17.22	178	212270	3.571	ng/ul#	86
15) Fluoranthene	19.16	202	2166465	25.876	ng/ul	84
17) Pyrene	19.52	202	2392993	28.165	ng/ul#	87
18) Benzo(a)anthracene	21.25	228	1209132	15.734	ng/ul#	84
19) Chrysene	21.30	228	1611736	20.218	ng/ul	98
21) Benzo(b)fluoranthene	22.95	252	2552232	35.131	ng/ul	88
22) Benzo(k)fluoranthene	22.99	252	846058m	10.965	ng/ul	
23) Benzo(a)pyrene	23.57	252	1160450m	16.894	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.19	276	713063m	8.965	ng/ul	
25) Dibenzo(a,h)anthracene	26.20	278	194744m	2.907	ng/ul	
26) Benzo(g,h,i)perylene	26.97	276	516456	7.572	ng/ul	97

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

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