

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061517\
 Data File : BE093231.D
 Acq On : 16 Jun 2017 5:07
 Operator : SJ/MA
 Sample : I3520-14DL 10X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5B41DL

Manual Integrations
 APPROVED

mohammad
 6/19/2017 11:07:30 AM

Quant Time: Jun 16 06:41:05 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 05:36:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2810	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	13890	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	7784	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	18108	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	19857	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	16504	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	681	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	1635	0.03	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	4444	0.129	ng/ul#	94
5) 2-Methylnaphthalene	12.18	142	972	0.041	ng/ul	98
7) Acenaphthylene	14.09	152	17827	0.534	ng/ul	96
8) Acenaphthene	14.43	153	663	0.025	ng/ul#	91
9) Fluorene	15.42	166	923	0.029	ng/ul#	83
12) Phenanthrene	17.14	178	10306	0.207	ng/ul#	89
13) Anthracene	17.22	178	39044	0.900	ng/ul#	88
15) Fluoranthene	19.16	202	83766	1.371	ng/ul	83
17) Pyrene	19.52	202	138002	2.357	ng/ul#	82
18) Benzo(a)anthracene	21.24	228	114610	2.164	ng/ul#	86
19) Chrysene	21.29	228	130409	2.373	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	327839m	6.546	ng/ul	
22) Benzo(k)fluoranthene	22.98	252	108489m	2.039	ng/ul	
23) Benzo(a)pyrene	23.56	252	143852	3.038	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.18	276	92483	1.687	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.20	278	24719	0.535	ng/ul	91
26) Benzo(g,h,i)perylene	26.96	276	68285	1.452	ng/ul	96

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

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