

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061417\
 Data File : BE093192.D
 Acq On : 15 Jun 2017 4:57
 Operator : SJ/MA
 Sample : I3521-05
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5C48

Manual Integrations
 APPROVED

mohammad
 6/15/2017 3:54:09 PM

Quant Time: Jun 15 06:35:56 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	3375	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	18081	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	11309	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	26425	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	27167	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	21972m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	5978	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	17001	0.24	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	5814	0.130	ng/ul#	94
5) 2-Methylnaphthalene	12.19	142	1269	0.041	ng/ul	96
7) Acenaphthylene	14.09	152	68782	1.417	ng/ul	95
8) Acenaphthene	14.43	153	1411	0.036	ng/ul#	93
9) Fluorene	15.42	166	1788	0.039	ng/ul#	88
12) Phenanthrene	17.13	178	40027	0.551	ng/ul#	85
13) Anthracene	17.24	178	59254	0.936	ng/ul	92
15) Fluoranthene	19.16	202	666101	7.471	ng/ul	84
17) Pyrene	19.52	202	803358	10.027	ng/ul#	84
18) Benzo(a)anthracene	21.24	228	402775	5.558	ng/ul#	83
19) Chrysene	21.30	228	641171	8.529	ng/ul	98
21) Benzo(b)fluoranthene	22.94	252	985551	14.781	ng/ul	86
22) Benzo(k)fluoranthene	22.98	252	352079m	4.972	ng/ul	
23) Benzo(a)pyrene	23.56	252	421233m	6.681	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.19	276	322857m	4.423	ng/ul	
25) Dibenzo(a,h)anthracene	26.20	278	84503m	1.375	ng/ul	
26) Benzo(g,h,i)perylene	26.97	276	254367	4.063	ng/ul	97

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

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