

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061517\
 Data File : BE093221.D
 Acq On : 15 Jun 2017 23:00
 Operator : SJ/MA
 Sample : I3522-16DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5B17DL

Manual Integrations
 APPROVED

mohammad
 6/19/2017 11:06:59 AM

Quant Time: Jun 16 05:31:19 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	3137	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	16021	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	9476	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	22737	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	24333	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	20669	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	1422	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	3807	0.06	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	8348	0.210	ng/ul#	92
5) 2-Methylnaphthalene	12.19	142	2065	0.076	ng/ul	98
7) Acenaphthylene	14.08	152	42958	1.056	ng/ul	96
8) Acenaphthene	14.43	153	1733	0.053	ng/ul	96
9) Fluorene	15.42	166	1861	0.048	ng/ul	89
12) Phenanthrene	17.14	178	34236	0.547	ng/ul#	87
13) Anthracene	17.22	178	68517	1.258	ng/ul#	86
15) Fluoranthene	19.16	202	414557	5.404	ng/ul	83
17) Pyrene	19.52	202	549408	7.656	ng/ul#	85
18) Benzo(a)anthracene	21.25	228	382323	5.890	ng/ul#	85
19) Chrysene	21.30	228	396881	5.895	ng/ul	98
21) Benzo(b)fluoranthene	22.94	252	800125m	12.756	ng/ul	
22) Benzo(k)fluoranthene	22.98	252	261540m	3.926	ng/ul	
23) Benzo(a)pyrene	23.56	252	360103	6.072	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.19	276	207759	3.025	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.19	278	61021	1.055	ng/ul	92
26) Benzo(g,h,i)perylene	26.96	276	144482	2.453	ng/ul	97

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

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