

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062017\
 Data File : BE093289.D
 Acq On : 20 Jun 2017 19:49
 Operator : SJ/JU
 Sample : I3627-18
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A91

Manual Integrations
 APPROVED

mohammad
 6/21/2017 12:18:36 PM

Quant Time: Jun 21 03:22:05 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 21 01:49:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1809	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	9364	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	6075	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	14119	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	14246	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	11471m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	3694	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	10146	0.26	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	80614	3.470	ng/ul#	90
5) 2-Methylnaphthalene	12.18	142	23975	1.509	ng/ul	98
7) Acenaphthylene	14.09	152	2415	0.093	ng/ul#	90
8) Acenaphthene	14.42	153	11563	0.550	ng/ul	97
9) Fluorene	15.42	166	28668	1.163	ng/ul	93
12) Phenanthrene	17.14	178	351158	9.039	ng/ul#	89
13) Anthracene	17.22	178	47482	1.404	ng/ul#	89
15) Fluoranthene	19.15	202	265180	5.566	ng/ul	84
17) Pyrene	19.52	202	167199	3.980	ng/ul#	84
18) Benzo(a)anthracene	21.24	228	46226	1.216	ng/ul#	85
19) Chrysene	21.29	228	36411	0.924	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	22054	0.634	ng/ul	90
22) Benzo(k)fluoranthene	22.97	252	8627m	0.233	ng/ul	
23) Benzo(a)pyrene	23.55	252	13709m	0.416	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.18	276	4760m	0.125	ng/ul	
25) Dibenzo(a,h)anthracene	26.19	278	1596m	0.050	ng/ul	
26) Benzo(g,h,i)perylene	26.96	276	3846	0.118	ng/ul#	89

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

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