

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061917\
 Data File : BE093272.D
 Acq On : 20 Jun 2017 4:24
 Operator : SJ/MA
 Sample : I3599-03
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5D37

Manual Integrations
 APPROVED

mohammad
 6/20/2017 3:13:38 PM

Quant Time: Jun 20 09:37:29 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 20 08:49:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2319	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	11748	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	7109	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	16657	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	17728	0.40	ng/ul	0.00
20) Perylene-d12	23.68	264	18980	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	2790	0.16	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	7509	0.17	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	77626	2.663	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	14505	0.728	ng/ul	99
7) Acenaphthylene	14.09	152	164104	5.379	ng/ul	95
8) Acenaphthene	14.42	153	14584	0.593	ng/ul	97
9) Fluorene	15.42	166	18231	0.632	ng/ul	93
12) Phenanthrene	17.14	178	212386	4.634	ng/ul#	88
13) Anthracene	17.22	178	314915	7.892	ng/ul#	87
15) Fluoranthene	19.16	202	1823420	32.443	ng/ul	83
17) Pyrene	19.52	202	2275431	43.523	ng/ul#	90
18) Benzo(a)anthracene	21.25	228	1175149	24.851	ng/ul#	84
19) Chrysene	21.30	228	1498925	30.557	ng/ul	97
21) Benzo(b)fluoranthene	22.96	252	2599053m	45.123	ng/ul	
22) Benzo(k)fluoranthene	22.99	252	921449m	15.063	ng/ul	
23) Benzo(a)pyrene	23.58	252	1210954	22.235	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.21	276	826069	13.099	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.21	278	246477	4.641	ng/ul	93
26) Benzo(g,h,i)perylene	26.99	276	510577	9.441	ng/ul	97

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

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