

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062017\
 Data File : BE093304.D
 Acq On : 21 Jun 2017 5:01
 Operator : SJ/JU
 Sample : I3599-03DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5D37DL

Manual Integrations
 APPROVED

mohammad
 6/21/2017 12:19:29 PM

Quant Time: Jun 21 07:05:15 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 03:23:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1878	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	10002	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5980	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	14716	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	15986	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	13984	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	535	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	1533	0.04	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	14800	0.596	ng/ul#	92
5) 2-Methylnaphthalene	12.18	142	2750	0.162	ng/ul	99
7) Acenaphthylene	14.09	152	33584	1.309	ng/ul	95
8) Acenaphthene	14.42	153	2917	0.141	ng/ul	96
9) Fluorene	15.42	166	3710	0.153	ng/ul	97
12) Phenanthrene	17.13	178	42805	1.057	ng/ul#	90
13) Anthracene	17.22	178	67671	1.920	ng/ul#	89
15) Fluoranthene	19.15	202	402573	8.107	ng/ul	83
17) Pyrene	19.51	202	519215	11.014	ng/ul#	87
18) Benzo(a)anthracene	21.24	228	257157	6.031	ng/ul#	85
19) Chrysene	21.30	228	341196	7.714	ng/ul	99
21) Benzo(b)fluoranthene	22.94	252	589662m	13.895	ng/ul	
22) Benzo(k)fluoranthene	22.98	252	206542m	4.583	ng/ul	
23) Benzo(a)pyrene	23.55	252	265990	6.629	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.18	276	173181	3.727	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.19	278	51520	1.317	ng/ul	92
26) Benzo(g,h,i)perylene	26.96	276	105172	2.640	ng/ul	96

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

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