

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062017\
 Data File : BE093282.D
 Acq On : 20 Jun 2017 15:29
 Operator : SJ/JU
 Sample : I3625-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A86

Manual Integrations
 APPROVED

mohammad
 6/21/2017 12:18:15 PM

Quant Time: Jun 21 02:20:23 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	1746	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	8839	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5671	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13404m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13849	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	11324m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	3830	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	10716	0.29	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.56	128	169792	7.742	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	23925	1.595	ng/ul	98
7) Acenaphthylene	14.09	152	2372	0.097	ng/ul#	94
8) Acenaphthene	14.42	153	26591	1.354	ng/ul	97
9) Fluorene	15.42	166	33832	1.470	ng/ul#	96
12) Phenanthrene	17.14	178	184936	5.014	ng/ul#	88
13) Anthracene	17.22	178	28255	0.880	ng/ul#	89
15) Fluoranthene	19.15	202	126915	2.806	ng/ul	84
17) Pyrene	19.51	202	91673	2.245	ng/ul#	87
18) Benzo(a)anthracene	21.24	228	23200	0.628	ng/ul#	85
19) Chrysene	21.30	228	19361	0.505	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	10904	0.317	ng/ul	90
22) Benzo(k)fluoranthene	22.97	252	3784m	0.104	ng/ul	
23) Benzo(a)pyrene	23.56	252	6425m	0.198	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.19	276	1672m	0.044	ng/ul	
25) Dibenzo(a,h)anthracene	26.19	278	599m	0.019	ng/ul	
26) Benzo(g,h,i)perylene	26.96	276	1189	0.037	ng/ul#	72

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

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