

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072517\
 Data File : BE093587.D
 Acq On : 25 Jul 2017 21:06
 Operator : SJ/JU
 Sample : I4209-20
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B24

Manual Integrations
 APPROVED

Sohil
 7/26/2017 5:56:04 PM

Quant Time: Jul 26 10:33:56 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 06:10:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	2279	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	12280	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	8238	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	20702m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	22883	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	20282m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	5154	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	17972	0.28	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.76	128	1180	0.040	ng/ul#	90
7) Acenaphthylene	14.25	152	1209	0.035	ng/ul#	87
12) Phenanthrene	17.29	178	1832	0.032	ng/ul#	91
13) Anthracene	17.38	178	1386	0.025	ng/ul#	55
15) Fluoranthene	19.30	202	16253	0.231	ng/ul	84
17) Pyrene	19.67	202	24415	0.366	ng/ul#	81
18) Benzo(a)anthracene	21.39	228	12855	0.194	ng/ul#	84
19) Chrysene	21.45	228	11027	0.173	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	16294	0.260	ng/ul	92
22) Benzo(k)fluoranthene	23.19	252	6209m	0.100	ng/ul	
23) Benzo(a)pyrene	23.81	252	15481m	0.259	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	8365m	0.127	ng/ul	
25) Dibenzo(a,h)anthracene	26.56	278	1817m	0.033	ng/ul	
26) Benzo(g,h,i)perylene	27.37	276	8917m	0.161	ng/ul	

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

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