

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072517\
 Data File : BE093581.D
 Acq On : 25 Jul 2017 17:25
 Operator : SJ/JU
 Sample : I4209-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B06

Manual Integrations
 APPROVED

Sohil
 7/26/2017 5:55:50 PM

Quant Time: Jul 25 18:49:00 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 18:00:47 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	2321	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	12463	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	8386	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	20425m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	22105	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	18237m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	6447	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	21080	0.33	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	2611	0.086	ng/ul#	96
8) Acenaphthene	14.59	153	1256	0.046	ng/ul	98
12) Phenanthrene	17.29	178	4906	0.088	ng/ul#	89
13) Anthracene	17.38	178	1976	0.036	ng/ul#	75
15) Fluoranthene	19.30	202	13570	0.195	ng/ul	84
17) Pyrene	19.67	202	17607	0.273	ng/ul#	82
18) Benzo(a)anthracene	21.39	228	12360	0.193	ng/ul#	88
19) Chrysene	21.45	228	12426	0.202	ng/ul	96
21) Benzo(b)fluoranthene	23.14	252	23647	0.419	ng/ul	91
22) Benzo(k)fluoranthene	23.19	252	8785m	0.158	ng/ul	
23) Benzo(a)pyrene	23.81	252	21608m	0.403	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	15267m	0.258	ng/ul	
25) Dibenzo(a,h)anthracene	26.56	278	4222m	0.085	ng/ul	
26) Benzo(g,h,i)perylene	27.36	276	15018m	0.301	ng/ul	

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

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