

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
 Data File : BE093567.D
 Acq On : 25 Jul 2017 8:49
 Operator : SJ/JU
 Sample : I4209-06
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B09

Manual Integrations
 APPROVED

Sohil
 7/25/2017 5:34:12 PM

Quant Time: Jul 25 12:09:00 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 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	2043	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	10840	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	7279	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	18590m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	22842	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	18946m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	5226	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	18417	0.31	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.76	128	1269	0.048	ng/ul#	89
12) Phenanthrene	17.29	178	1808	0.036	ng/ul#	90
15) Fluoranthene	19.30	202	5042	0.080	ng/ul	85
17) Pyrene	19.67	202	4974	0.075	ng/ul#	87
18) Benzo(a)anthracene	21.40	228	3527	0.053	ng/ul#	88
19) Chrysene	21.45	228	3690m	0.058	ng/ul	
21) Benzo(b)fluoranthene	23.15	252	5844	0.100	ng/ul	84
22) Benzo(k)fluoranthene	23.19	252	2451m	0.042	ng/ul	
23) Benzo(a)pyrene	23.80	252	4607m	0.083	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	3144m	0.051	ng/ul	
26) Benzo(g,h,i)perylene	27.37	276	3096m	0.060	ng/ul	

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

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