

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
 Data File : BE093549.D
 Acq On : 24 Jul 2017 21:47
 Operator : SJ/JU
 Sample : I4213-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B29

Manual Integrations
 APPROVED

Sohil
 7/25/2017 5:33:11 PM

Quant Time: Jul 25 04:23:40 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 24 18:07:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	3771	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	20323	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	13284	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	30586m	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	34550	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	30829m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	8761	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	29286	0.30	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.76	128	2309	0.047	ng/ul#	93
7) Acenaphthylene	14.25	152	3019	0.054	ng/ul	97
9) Fluorene	15.57	166	1476	0.027	ng/ul	92
12) Phenanthrene	17.29	178	18479m	0.222	ng/ul	
13) Anthracene	17.40	178	10161	0.125	ng/ul	94
15) Fluoranthene	19.31	202	73276	0.703	ng/ul	83
17) Pyrene	19.67	202	63060	0.625	ng/ul#	81
18) Benzo(a)anthracene	21.40	228	56658	0.566	ng/ul#	87
19) Chrysene	21.45	228	76574	0.797	ng/ul	99
21) Benzo(b)fluoranthene	23.15	252	58557	0.614	ng/ul	88
22) Benzo(k)fluoranthene	23.20	252	24496m	0.260	ng/ul	
23) Benzo(a)pyrene	23.81	252	38947m	0.429	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	19348m	0.193	ng/ul	
25) Dibenzo(a,h)anthracene	26.57	278	5266m	0.063	ng/ul	
26) Benzo(g,h,i)perylene	27.38	276	14955m	0.177	ng/ul	

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

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