

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
 Data File : BE093566.D
 Acq On : 25 Jul 2017 8:13
 Operator : SJ/JU
 Sample : I4209-05
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B08

Manual Integrations
 APPROVED

Sohil
 7/25/2017 5:34:10 PM

Quant Time: Jul 25 12:03:48 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	2260	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	12103	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	8019	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	20513m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	23046	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	19174m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4631	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	17302	0.27	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.76	128	1317	0.045	ng/ul#	89
12) Phenanthrene	17.29	178	4013	0.072	ng/ul#	89
15) Fluoranthene	19.31	202	11319	0.162	ng/ul	84
17) Pyrene	19.67	202	14116	0.210	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	10607	0.159	ng/ul#	91
19) Chrysene	21.45	228	11188	0.175	ng/ul#	93
21) Benzo(b)fluoranthene	23.15	252	17049	0.287	ng/ul	93
22) Benzo(k)fluoranthene	23.19	252	6685m	0.114	ng/ul	
23) Benzo(a)pyrene	23.81	252	14352m	0.254	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	8858m	0.142	ng/ul	
26) Benzo(g,h,i)perylene	27.36	276	8085m	0.154	ng/ul	

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

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