

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081417\
 Data File : BE093795.D
 Acq On : 14 Aug 2017 20:06
 Operator : SJ/JU
 Sample : I4557-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC6

Manual Integrations
 APPROVED

Sohil
 8/15/2017 6:43:32 PM

Quant Time: Aug 15 11:04:02 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 Aug 15 10:51:12 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2562	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	13464	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	8948	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	18265	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	18407	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	15761m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	6831	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	18479	0.32	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.75	128	13318	0.407	ng/ul#	95
5) 2-Methylnaphthalene	12.36	142	23191	0.971	ng/ul	99
7) Acenaphthylene	14.24	152	6281	0.168	ng/ul#	73
8) Acenaphthene	14.58	153	9144	0.311	ng/ul	95
9) Fluorene	15.57	166	34993	0.940	ng/ul	91
12) Phenanthrene	17.29	178	112791	2.267	ng/ul#	90
13) Anthracene	17.38	178	35060	0.724	ng/ul#	86
15) Fluoranthene	19.30	202	143957	2.314	ng/ul	83
17) Pyrene	19.67	202	173703	3.233	ng/ul#	81
18) Benzo(a)anthracene	21.40	228	71921	1.348	ng/ul#	91
19) Chrysene	21.45	228	66584	1.301	ng/ul#	90
21) Benzo(b)fluoranthene	23.15	252	68175m	1.398	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	23565m	0.490	ng/ul	
23) Benzo(a)pyrene	23.81	252	53214m	1.147	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	28367m	0.554	ng/ul	
25) Dibenzo(a,h)anthracene	26.57	278	8048m	0.188	ng/ul	
26) Benzo(g,h,i)perylene	27.37	276	23253m	0.539	ng/ul	

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

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