

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102617\
 Data File : BE094565.D
 Acq On : 26 Oct 2017 16:49
 Operator : SJ/JU
 Sample : I5873-05 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-16(7-7.75) 1012117

Manual Integrations
 APPROVED

Sohil
 10/27/2017 5:34:53 PM

Quant Time: Oct 27 08:28:00 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 27 04:05:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1305	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	7728	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.52	164	4081m	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8339	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	8366	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	9264	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	139626	11.13	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	82919	3.35	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	246885	13.066	ng/ul#	94
5) 2-Methylnaphthalene	12.41	142	310807	23.512	ng/ul	98
7) Acenaphthylene	14.24	152	8847m	0.479	ng/ul	
8) Acenaphthene	14.59	153	23321	1.758	ng/ul#	68
9) Fluorene	15.57	166	43197	2.900	ng/ul#	85
12) Phenanthrene	17.29	178	293152	13.309	ng/ul#	89
13) Anthracene	17.38	178	14139	0.650	ng/ul#	74
15) Fluoranthene	19.31	202	16237	0.638	ng/ul	79
17) Pyrene	19.67	202	46375	1.793	ng/ul#	88
18) Benzo(a)anthracene	21.40	228	8694	0.364	ng/ul#	47
19) Chrysene	21.46	228	14523	0.587	ng/ul#	63
21) Benzo(b)fluoranthene	23.17	252	5330m	0.205	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	1779m	0.064	ng/ul	
23) Benzo(a)pyrene	23.84	252	6678	0.256	ng/ul#	44
24) Indeno(1,2,3-cd)pyrene	26.64	276	2309	0.086	ng/ul#	88
25) Dibenzo(a,h)anthracene	26.62	278	1266	0.056	ng/ul#	1
26) Benzo(g,h,i)perylene	27.48	276	3277	0.151	ng/ul#	44

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

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