

Data Path : Z:\HPCHEM1\BNA E\DATA\BE103117\
 Data File : BE094669.D
 Acq On : 31 Oct 2017 21:50
 Operator : SJ/JU
 Sample : I5873-05RE 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-16(7-7.75)_1012117RE

Manual Integrations
 APPROVED

Sohil
 11/1/2017 4:29:46 PM

Quant Time: Nov 01 04:33:29 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 02:55:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1284	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	7776	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.52	164	4169m	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8728	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	8256	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	9481	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	159791	12.66	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	94039	3.63	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.74	128	263875	13.879	ng/ul#	93
5) 2-Methylnaphthalene	12.41	142	338441	25.445	ng/ul	98
7) Acenaphthylene	14.24	152	9832m	0.521	ng/ul	
8) Acenaphthene	14.59	153	26263	1.938	ng/ul#	70
9) Fluorene	15.57	166	47106	3.096	ng/ul#	87
12) Phenanthrene	17.29	178	318496	13.816	ng/ul#	89
13) Anthracene	17.38	178	15490	0.681	ng/ul#	75
15) Fluoranthene	19.31	202	17484	0.656	ng/ul	79
17) Pyrene	19.67	202	49670	1.946	ng/ul#	87
18) Benzo(a)anthracene	21.41	228	9668	0.410	ng/ul#	46
19) Chrysene	21.46	228	15229	0.624	ng/ul#	63
21) Benzo(b)fluoranthene	23.18	252	5766m	0.216	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	2012m	0.070	ng/ul	
23) Benzo(a)pyrene	23.85	252	6941	0.260	ng/ul#	19
24) Indeno(1,2,3-cd)pyrene	26.66	276	2070	0.075	ng/ul#	47
25) Dibenzo(a,h)anthracene	26.64	278	1143	0.049	ng/ul#	1
26) Benzo(a,h,i)perylene	27.51	276	2897	0.131	ng/ul#	9

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

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