

Data Path : Z:\HPCHEM1\BNA E\DATA\BE110515\
 Data File : BE091074.D
 Acq On : 5 Nov 2015 19:40
 Operator : UM/NP
 Sample : G4273-11
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AP7

Manual Integrations
 APPROVED

Mmdadoda
 11/6/2015 3:45:22 PM

Quant Time: Nov 06 06:40:52 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM02.2-EPA-SIM-BE102715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 06 05:41:57 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	3552	0.40	ng/ul	0.00
4) Naphthalene-d8	9.57	136	20110	0.40	ng/ul	0.00
8) Acenaphthene-d10	13.40	164	11977	0.40	ng/ul	0.00
12) Phenanthrene-d10	16.14	188	28963m	0.40	ng/ul	0.00
18) Chrysene-d12	20.29	240	31405	0.40	ng/ul	0.00
22) Perylene-d12	22.19	264	41055m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	2.93	96	14824	5.80	ng/uL	-0.02
6) 2-Methylnaphthalene-d10	11.07	152	5543	0.20	ng/ul	0.00
16) Fluoranthene-d10	18.13	212	17310	0.21	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	9.62	128	26576	0.54	ng/ul	99
7) 2-Methylnaphthalene	11.14	142	50110	1.55	ng/ul	99
9) Acenaphthylene	13.14	152	18673	0.31	ng/ul	99
10) Acenaphthene	13.46	153	188159	4.59	ng/ul	94
11) Fluorene	14.45	166	230715	4.39	ng/ul	89
14) Phenanthrene	16.17	178	4381400m	13.10	ng/ul	
15) Anthracene	16.25	178	1783991	5.87	ng/ul	97
17) Fluoranthene	18.15	202	3798063m	9.70	ng/ul	
19) Pyrene	18.54	202	2701721	7.57	ng/ul#	78
20) Benzo(a)anthracene	20.26	228	295700	3.67	ng/ul	96
21) Chrysene	20.32	228	270904	2.92	ng/ul	98
23) Benzo(b)fluoranthene	21.63	252	161429	1.26	ng/ul	94
24) Benzo(k)fluoranthene	21.65	252	181428m	1.38	ng/ul	
25) Benzo(a)pyrene	22.10	252	196481m	1.66	ng/ul	
26) Indeno(1,2,3-cd)pyrene	23.83	276	393230m	0.79	ng/ul	
27) Dibenzo(a,h)anthracene	23.79	278	27598m	0.23	ng/ul	
28) Benzo(g,h,i)perylene	24.40	276	317216m	0.61	ng/ul	

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

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