

Data Path : V:\HPCHEM1\BNA M\DATA\BM112115\
 Data File : BM003275.D
 Acq On : 22 Nov 2015 10:55
 Operator : UM/IZ
 Sample : G4345-12DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4J54DL

Manual Integrations
 APPROVED

Mmdadoda
 11/23/2015 7:05:11 PM

Quant Time: Nov 23 01:16:51 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM111815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 23 01:07:06 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	13262	0.40	ng/ul	0.00
4) Naphthalene-d8	10.52	136	53372	0.40	ng/ul	-0.01
8) Acenaphthene-d10	14.38	164	26192	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.12	188	50357	0.40	ng/ul	0.00
18) Chrysene-d12	21.33	240	43974	0.40	ng/ul	0.01
22) Perylene-d12	23.58	264	44029	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.20	96	8435	0.60	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.12	152	2655	0.04	ng/ul	0.00
16) Fluoranthene-d10	19.17	212	3993	0.03	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.57	128	3615	0.03	ng/ul#	89
7) 2-Methylnaphthalene	12.19	142	1673	0.02	ng/ul#	100
9) Acenaphthylene	14.09	152	4408	0.04	ng/ul#	93
13) Pentachlorophenol	16.78	266	685	0.06	ng/ul	94
14) Phenanthrene	17.16	178	70868m	0.48	ng/ul	
15) Anthracene	17.26	178	12024	0.10	ng/ul	96
17) Fluoranthene	19.19	202	143770	0.89	ng/ul	99
19) Pyrene	19.55	202	131987	0.88	ng/ul	99
20) Benzo(a)anthracene	21.31	228	58741	0.46	ng/ul	95
21) Chrysene	21.36	228	63404	0.44	ng/ul	98
23) Benzo(b)fluoranthene	22.91	252	86581m	0.52	ng/ul	
24) Benzo(k)fluoranthene	22.93	252	32724m	0.21	ng/ul	
25) Benzo(a)pyrene	23.48	252	58342	0.39	ng/ul	98
26) Indeno(1,2,3-cd)pyrene	25.86	276	42683	0.24	ng/ul#	94
27) Dibenzo(a,h)anthracene	25.87	278	9728	0.07	ng/ul#	73
28) Benzo(a,h,i)perylene	26.57	276	43319	0.28	ng/ul	98

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

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