

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122216\
 Data File : BM008574.D
 Acq On : 23 Dec 2016 02:16
 Operator : UM/SJ
 Sample : H5970-17DL 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA856DL

Manual Integrations
 APPROVED

umangi
 12/23/2016 6:12:02 PM

Quant Time: Dec 23 05:57:16 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 23 05:50:04 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	369	0.40	ng/ul	-0.06
2) Naphthalene-d8	10.45	136	1140	0.40	ng/ul	-0.04
6) Acenaphthene-d10	14.30	164	665	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.06	188	1179	0.40	ng/ul	-0.02
16) Chrysene-d12	21.25	240	1256	0.40	ng/ul	-0.02
20) Perylene-d12	23.48	264	1417	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	84	0.05	ng/ul	-0.02
14) Fluoranthene-d10	19.10	212	231	0.07	ng/ul	-0.01
Target Compounds						
					Ovalue	
3) Naphthalene	10.50	128	435	0.14	ng/ul#	52
5) 2-Methylnaphthalene	12.14	142	422	0.20	ng/ul	95
7) Acenaphthylene	14.01	152	903	0.30	ng/ul#	85
8) Acenaphthene	14.35	153	156	0.07	ng/ul#	81
9) Fluorene	15.36	166	186	0.07	ng/ul#	74
12) Phenanthrene	17.09	178	2135	0.59	ng/ul	98
13) Anthracene	17.19	178	1371m	0.39	ng/ul	
15) Fluoranthene	19.12	202	4902	1.11	ng/ul	96
17) Pyrene	19.48	202	8231	1.70	ng/ul	97
18) Benzo(a)anthracene	21.24	228	5566m	1.35	ng/ul	
19) Chrysene	21.28	228	6923m	1.53	ng/ul	
21) Benzo(b)fluoranthene	22.81	252	14084m	2.52	ng/ul	
22) Benzo(k)fluoranthene	22.85	252	4298m	0.79	ng/ul	
23) Benzo(a)pyrene	23.38	252	7233m	1.40	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.71	276	5451	0.88	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.71	278	1465	0.30	ng/ul#	87
26) Benzo(a,h,i)perylene	26.40	276	5030	0.95	ng/ul#	92

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

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