

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122716\
 Data File : BM008679.D
 Acq On : 27 Dec 2016 18:37
 Operator : UM/SJ
 Sample : H6069-01DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA7H7DL

Manual Integrations
 APPROVED

sohil
 12/28/2016 9:50:27 AM

Quant Time: Dec 28 04:50:07 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 28 03:24:04 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	351m	0.40	ng/ul	-0.17
2) Naphthalene-d8	10.45	136	1198	0.40	ng/ul	-0.12
6) Acenaphthene-d10	14.30	164	765m	0.40	ng/ul	-0.05
10) Phenanthrene-d10	17.06	188	1351	0.40	ng/ul	-0.07
16) Chrysene-d12	21.26	240	1639	0.40	ng/ul	-0.03
20) Perylene-d12	23.49	264	1548	0.40	ng/ul	-0.04
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.08	152	73m	0.04	ng/ul	-0.09
14) Fluoranthene-d10	19.10	212	222m	0.06	ng/ul	-0.05
Target Compounds						
15) Fluoranthene	19.12	202	985	0.19	ng/ul	67
17) Pyrene	19.48	202	1117	0.18	ng/ul#	75
18) Benzo(a)anthracene	21.25	228	1004	0.19	ng/ul#	28
19) Chrysene	21.30	228	575	0.10	ng/ul#	23
21) Benzo(b)fluoranthene	22.83	252	1605m	0.26	ng/ul	
22) Benzo(k)fluoranthene	22.85	252	650m	0.11	ng/ul	
23) Benzo(a)pyrene	23.39	252	894	0.16	ng/ul#	22
24) Indeno(1,2,3-cd)pyrene	25.76	276	811	0.12	ng/ul	97
26) Benzo(g,h,i)perylene	26.44	276	732	0.13	ng/ul#	31

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

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